EPA Region 5 Records Ctr.

SITE ASSESSMENT REPORT

FOR

STANDARD SCRAP PROCESSING

(aka: SCRAP METAL)

U.S. EPA ID: ILD045698263

SS ID: NA

TDD: T05-9402-007 PAN: EIL0831SAA

MAY 6, 1994

Reviewed by

Date: 5-6-94

Date: 5/6/94

Approved by



ecology and environment, inc.

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recycled paper

NOTICE

Originally, this site was tasked to Ecology & Environment, Inc. under the name Scrap Metal. The actual name of this site is Standard Scrap Processing. All file information, photologs, maps, and text in this report refer to the site by its original tasked name Scrap Metal.

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1.0 INTRODUCTION

The Ecology & Environment, Inc. (E & E), Technical Assistance Team (TAT) was tasked by the United States Environmental Protection Agency (U.S. EPA) to perform a site assessment at the Scrap Metal (SM) site under Technical Directive Document (TDD) T05-9402-007. Activities performed for the site assessment included prepared and implemented Health and Safety Plan; compiled available information; prepared and implemented sampling plan; evaluated threat to human health and environment; and provided photodocumentation. The following report provides a summary of these activities.

2.0 SITE DESCRIPTION

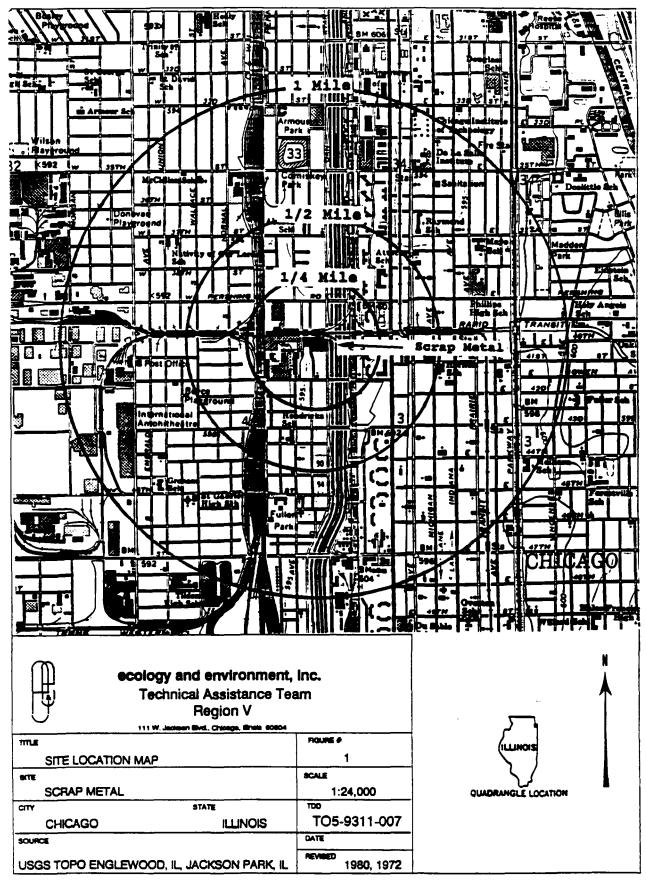
The SM site is located at 4004 South Wentworth Avenue, Chicago, Cook County, Illinois, Latitude 87° 37' 55" north, Longitude 41° 52' 50" west, (see Figure 1, Site Location Map). The SM site is located in a industrial and residential area. The facility is an active 3-acre scrap yard involved in the reclamation of metals. Past and present operations have taken place on two distinct parcels of property separated by Wells Avenue.

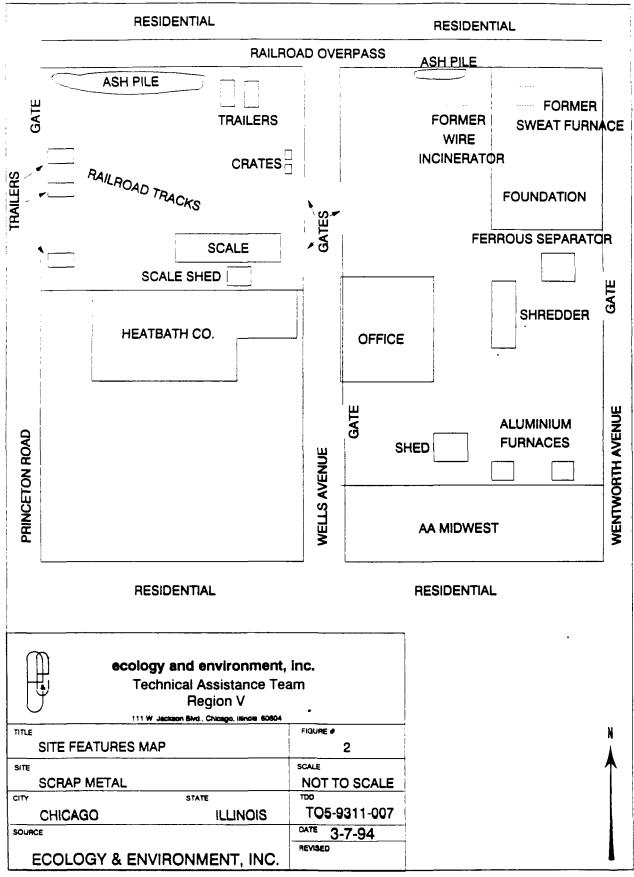
The east lot is approximately 2.5 acres in size and the west lot is approximately 0.5 acres (see Figure 2, Site Features map). The east lot has a office building, a small receiving shed, two aluminum furnaces, a ferrous metal sorter, a shredder, and large piles of scrap metal located at various points throughout the property. The west lot has a small scale house used to weigh incoming scrap metal trucks and is also used to park semitrailers. A railroad spur cuts across both lots diagonally.

3.0 SITE BACKGROUND

The property has been used for industrial purposes since 1895. A Sanborn Fire Insurance map, dated, 1895 indicates part of the east and west lots was owned by Weaver and Getz. It is unknown what business Weaver and Getz operated. Two other businesses owned parts of the east lot; W. B. Scace and Company (which loaded lime and cement) and Chas D. Pek and Company (which sold hay, grain, and feed). A Sanborn Fire Insurance map, dated, 1925, shows a Baker-Smith Coal Company operated a coal yard on the east lot. The west lot did not appear to be used by either Chicago Mfg. & Distribution Company (a machine shop) or Arthur M. Adler & Company, Oils located to the south.

The Standard Scrap Metal Company (SSMC) was started in 1928 by Sam Cohen and Sam Kanter at 4004 Wentworth Avenue. SSMC was involved in reclaiming aluminum and cooper and sold the reclaimed scrap metal to steel smelters and refiners. The facility contained 1 gas-fired boiler, 2 aluminum sweat furnaces, and a





wire burning incinerator. Operations continued until 1972 when the company merged into Standard Scrap Metal Company, Incorporated (SSCMI). SSCMI continued operations at the site until 1987, when the company went bankrupt.

Phoenix Recycling, a metals reclamation business, began operations at the site about the same time of the SSCMI bankruptcy and continued operating the same type of business until 1989. The Sam Cohen and Sam and Benjamin Kanter Building Partnership owned Phoenix Recycling.

Chicago International Exporting Company (CIEC) began operations at the site in 1989 and continues to operate a metal recycling facility. CIEC is owned by Chicago International Export Company, Incorporated whose President is Steve Cohen.

The SM site has been investigated by the Illinois Environmental Protection Agency (IEPA) and U.S. EPA beginning in 1973. In 1973, personnel from IEPA inspected the site for compliance with Air Pollution Regulations. The inspection revealed that the facility did not have the proper air pollution permits to operate their incinerator or sweat furnaces. A suit (PCB 83-22) was filed against SSMCI for not possessing permits required by IEPA and the City of Chicago. The complaint stated that SSMCI could achieve compliance by installing afterburners on the sweat furnaces. The afterburners were not installed and permits were not applied for until 1984. A permit for the gas-fired boiler was applied for and approved on December 14, 1984.

On January 10, 1985, Illinois Pollution Board (IPB) continued the suit (PCB 83-22) against SSMCI for permit violations. IPB suit ordered SSMCI to:

Cease and desist from operations of its incinerator until the necessary operating permit is obtained from the IEPA: Cease and desist from operating either of its aluminum sweat furnaces until the necessary permits are obtained from the IEPA and permanently shut down the inactive aluminum sweat furnace by January 21, 1985.

Install temperature gauges on each afterburner with an interlock that prevents operation unless the afterburner temperature is at least 1400 degrees Fahrenheit, and take all necessary steps to ensure adequate pre-heating of each afterburner prior to charging. These requirements are to be made conditions of the operating permits issued by the IEPA;

Within 90 days of the date of this order pay a penalty of \$30,000 for the violation of the Act and Regulations as described in this opinion.

On February 14, 1984, IEPA investigated a report from an employee of Heatbath Corporation, the plant south of the west lot, that workers at the facility periodically dumped transformer oil on the ground and igniting it. This practice was to have taken place from 1977 to 1981. On one occasion the roof of the Heatbath Corporation caught fire and the Chicago Fire Department extinguished the fire.

IEPA collected 2 soil samples, one sample from the west lot and one from near a garage located at 3949 South Wells Avenue. A complaint from the resident at 3949 South Wells Avenue stated oil from SM Site would flow off-site into her yard. Sampling results indicated Polychlorinated biphenyls (PCBs) at 3.9 parts per million (ppm) at the residence and 1,300 ppm from the west lot. The IEPA requested that the U.S. EPA conduct a PCB inspection at the site.

On March 30, 1984, U.S. EPA's Toxic Substance Office conducted a inspection of SM facility to document their handling, storage, and disposal practices. During the inspection 6 composite soil samples and 1 wipe sample were collected from the west lot and a residence at 3949 South Wells Avenue. Analytical results from the samples indicated PCB levels of up to 2,095 ppm but no detectable amounts at the residence. The U.S. EPA filed a complaint against SSMCI for violating regulations pertaining to the disposal of PCBs and a civil penalty of \$25,000 was levied for improper disposal of PCBs.

On June 18, 1985, the Roy F. Weston TAT collected 4 soil samples and 2 wipe samples (east lot) at the SM site. The analytical results indicated PCBs levels up to 336 ppm in three samples and some dioxin isomers in 4 samples.

On October 29, 1985, an amended complaint by U.S. EPA was filed against SSMCI after the June 18, 1985 inspection. The amended complaint levied a \$30,000 penalty for violations of Section 16(a) of the Toxic Substance Control Act (TSCA). In February 1987, SSMCI appealed the decision and the complaint was dismissed because the U.S. EPA did not prove that the PCB-contaminated oil had been accepted at the site after 1978. However, the U.S. EPA appealed the dismissal and the decision was reversed and the \$30,000 fine was levied against the facility. SSMCI filed for bankruptcy, the fine was not collected and the U.S. EPA Enforcement Branch closed the case. In 1989, metal recycling operations continued on-site, when CIEC began operation at the site.

In 1990, a former railroad employee had a telephone interview with Tom Crause of the IEPA. The former railroad employee indicated that workers at the facility cut up and disposed of many electrical transformers during his 30 years of working for the railroad. Based on this information, on August 27, 1990,

SSMC was placed on the Comprehensive Environmental Response, Compensation and Liability Information System (CERCLIS).

On August 29, 1991, IEPA personnel conducted an off-site reconnaissance inspection of the SM facility. IEPA observed piles of scrap metal around the site. No air admission were observed at the site, the boiler appeared to be not in operation. At the east lot, the north sweat furnace had been demolished and was left as a pile of debris. A number of drums, which appeared to be empty, were observed near the north side of the office building. No leakage was observed from the drums or stressed vegetation on the lot. At the west lot, the gates were open and was empty except for 3 semi-trailers. The IEPA prepared a Preliminary Assessment (PA) for Standard Scrap Metal site (also called Standard Metal) on September 30, 1991.

On September 22, 1992, the IEPA was tasked by the U.S. EPA Region V to conduct a CERCLIS Screening Site Inspection (SSI) of the SM site. The SSI was conducted on November 4 and 5, 1992 and consisted of the collection of 12 soil samples. The analytical results from on-site soil sampling indicated PCBs up to 670,000 Micrograms per kilogram (ug/kg = parts per billion (ppb). The PCBs can be directly associated with past activities at the site as reported by the Heatbath employee and railroad employee.

4.0 SITE ASSESSMENT

On February 22, 1994, TAT members John Nordine and Yvette Anderson met with U.S. EPA On-Scene-Coordinator (OSC) Steve Faryan at the SM site. The SM site is currently called Chicago International Export Company. Access to the site was granted during an interview with CIEC President, Steve Cohn and Secretary Treasure, Buddy Cohn. A site safety meeting was held and the potential hazards of the site were discussed. A site reconnaissance of the SM Site was completed, which included a tour of the facility (see Appendix A for Site Photo Log).

A railroad spur cuts across both lots diagonally. A railroad overpass delineates the north property line for both lots (see Figure 2, Site Features map). A 8-foot fence topped with 3 strands of barbed wire surround both lots.

The east lot has an office building, a small receiving shed, 2 aluminum furnaces, a ferrous metal sorter, a shredder, a ash pile, and large piles of scrap metal located at various points throughout the property. According to Buddy Cohn, the entire east lot has been paved with either concrete or asphalt. The TAT could not verify that the east lot had been paved due to debris and piles of scrap metal covering the lot. A debris and ash pile was located in the area where the wire incinerator was formerly located. The sweat furnaces have been demolished. A foundation

of a building was located in the northeast corner of the east lot.

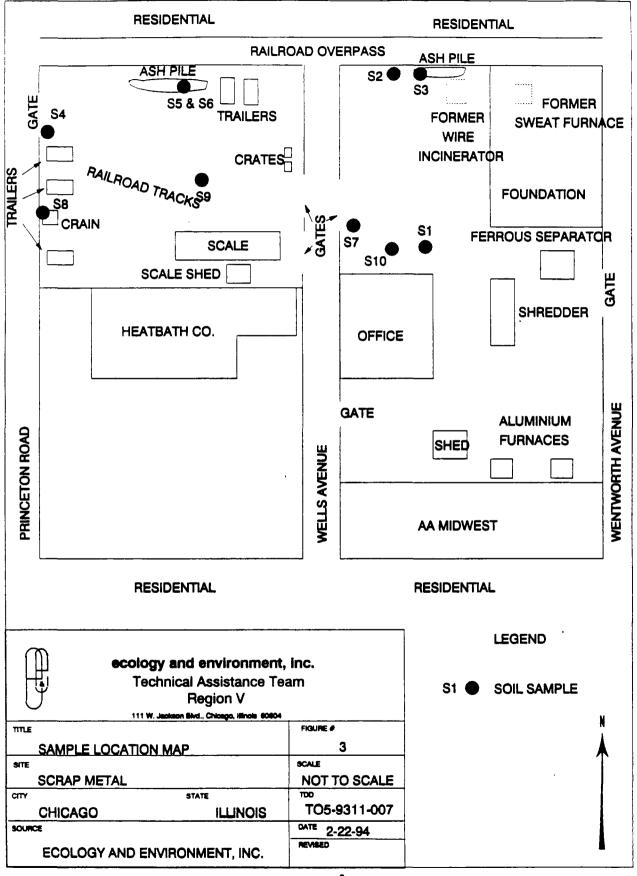
A small scale house is located on the west lot along the south side which is used to weigh incoming trucks loaded with scrap metal. A crane and 3 semi-trailers were parked along the west fence line of the west lot. Located along the north property line from west to east were a crane boom, a ash pile with tires placed on top, a striped out car, and 2 semi-trailers. Three large crates were located along the east fence line of the west lot.

During the reconnaissance CIEC workers were observed to burn a pile of cardboard material at the east lot and burn unknown materials in a 55-gallon drum at the west lot. Later CIEC workers put out the cardboard fire with water from a hose. The materials burned in the 55-gallon drum gave off a black smoke that was irritating to the eyes, nose, and throat. When the metal shredder was in operation a dust cloud could be observed. The metal shredder was turned off while TAT was in the area. Oil stains were observed on the ground at both lots. A motor had been cut open and oil was observed flowing from it at the east lot.

To characterize the possible hazardous substances on-site, the TAT collected ten soil samples including a duplicate sample and a matrix spike/matrix spike duplicate (MS/MSD) sample (see Figure 3 Sample Location Map). All soil samples were collected from 1 to 3 inches in depth with a shovel and a trowel.

All soil samples collected on-site were collected in Level D personal protection. Sampling gloves were changed before each sample was collected. Sampling equipment was decontaminated after each use, using a alconox and distilled water wash solution and triple rinsed with distilled water. All personal protective equipment was decontaminated as above and disposed of properly.

Soil samples S1, S2, S3, S7, and S10 were collected from the east lot. Soil sample S1 was collected from an area approximately 15 north of the northeast corner of the office. Soil Sample S2 was collected near a tree along the north property line west of the former wire incinerator. S3 was collected from a ash pile approximately 20 feet east of sample S2. An white crystalline substance was noted in the ash material and two dead rats were noted on top of the ash pile. Sample S7 was collected 25 feet north and 25 feet east of the northwest corner of the office building near the railroad tracks. Soil sample S10 was collected approximately 10 feet north and 35 feet east of the northwest corner of the office building.



Soil samples S4, S5, S6, S8, and S9 were collected from the west lot. Soil sample S4 was collected 3 feet west of the northwest gate next to the railroad tracks. S4 was also the Matrix Spike/Matrix Spike Duplicate (MS/MSD). Soil sample S5 and S6, a duplicate sample, were collected from the ash pile along the north property line. S8 was collected by the crane 30 feet north and 3 feet east of the northwest corner of the Heatbath Company building. Soil sample S9 was collected approximately from the middle of the west lot by the railroad tracks.

Soil samples were delivered to and analyzed by Twin City Testing Corporation, 737 Pelham Boulevard, St. Paul, Minnesota 55114. The soil samples were analyzed for semi volatile organics, total Resources Conservation Recovery Act (RCRA) 10 metals, Toxicity Characteristic Leaching Procedure (TCLP) for 8 metals, PCBs, and dioxin (soil samples S1, S2, S3, and S8).

5.0 ANALYTICAL RESULTS

Results of the TAT collected samples revealed the presence of U. S. EPA Target Compound List (TCL) volatile compounds and base neutral acids; Total RCRA 10 metals; Toxicity Characteristic Leaching Procedure (TCLP) 8 metals; PCBs; and Dioxin. See Appendix B for data quality assurance review. See Tables 1 through 4 for analytical results.

6.0 DISCUSSION OF POTENTIAL THREATS

Conditions at the Scrap Metal Site that may warrant a removal action, as set forth in paragraph (b) (2) of Section 300.415 of the National Contingency Plan (NCP), include:

- 1. Actual or potential exposure to nearby human populations, animals, or the food chain from hazardous substances or pollutants or contaminants; and
- 2. high levels of hazardous substances or pollutants or contaminants in soils largely at or near the surface, that may migrate.

6.1 Potential Exposure To Hazardous Substances

Site investigations and analytical data obtained by IEPA and the TAT have identified soil containing materials that possess characteristics of hazardous wastes, as defined by RCRA at the SM site. TCLP metals, cadmium at 1.3 mg/l and lead at 71 mg/l, were detected in on-site soils. On-site soils contain heavy metals that were above the RCRA limits for cadmium and lead (1 and 5 mg/L, respectively) which are listed waste D006 and D008, respectively. Total PCBs were detected in on-site soils ranging from 37,100 to 2,000,000

TABLE 1 ANALYTICAL DATA FOR TOTAL METALS AT THE SCRAP METAL SITE CHICAGO, ILLINOIS units = mg/Kg

SAMPLE NUMBER PARAMETER	S1	S2	S3	S4	S4*	S5	S6**	S7	S8	S9	S10
ARSENIC	36	56	22	36	39	35	34	43	38	13	50
BARIUM	410	1,400	450	160	130	540	550	630	390	190	1,300
CADMIUM	49	140	68	19	20	64	57	42	69	20	47
CHROMIUM	97	200	130	300	350	1,400	1,700	81	290	48	92
COPPER	20,000	15,000	38,000	1,600	1,600	15,000	8,500	7,400	10,000	7,400	6,400
LEAD	2,600	5,900	5,400	870	900	5,700	3,500	2,000	2,400	3,200	5,700
MERCURY	16	25	3.3	5.9	6.1	10	13	12	20	7.1	5.2
SELENIUM	21	82	ND	ND	5.5	10	11	13	ND	6.5	ND
SILVER	13	21	67	2.1	1.9	39	23	9.7	4.3	9.6	2.2
ZINC	9,000	6,500	3,600	1,200	1,100	5,400	4,600	4,800	4,200	2,600	9,400

^{* =} Sample S4 is a Matrix Spike/Matrix Spike Duplicate sample.

ND = Non detect

NOTE: Correction made to units (from ppb to mg/Kg) per laboratory data sheets (12/2/94)

^{** =} Sample S6 is a duplicate sample of S5.

TABLE 2
ANALYTICAL DATA FOR TCLP METALS
AT THE SCRAP METAL SITE
CHICAGO, ILLINOIS
units = mg/l

SAMPLE NUMBER PARAMETER	S1	S2	S3	S4	S4*	S5	S6**	S7	S8	S9	S10	REGUL- ATORY LEVEL
ARSENIC	ND	ND	ND.	ND	ND	ND	ND	ND	ИD	ND	ND	5.0
BARIUM	1.5	2.8	2.7	1.1	1.1	4.8	2.7	1.3	1.5	1.6	2.0	100
CADMIUM	0.87	1.3	0.11	0.23	0.22	0.91	0.66	1.2	0.85	0.48	0.69	1.0
CHROMIUM	ND	ND	ND	0.01	ND	ND	ND	ND	0.012	0.015	ND	5.0
LEAD	7.5	16	0.69	8.2	8.1	45	4.3	7.4	5.3	71	32-	5.0
MERCURY	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.2
SELENIUM	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.0
SILVER	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	5.0

^{* =} Sample S4 is a Matrix Spike/Matrix Spike Duplicate sample.

ND = Non detected

NOTE: Corrections made to metal values that are in bold per laboratory data sheets (12/2/94)

^{** =} Sample S6 is a duplicate sample of S5.

TABLE 3
ANALYTICAL DATA FOR PCBs
AT THE SCRAP METAL SITE
CHICAGO, ILLINOIS
unit = ppb

SAMPLE NUMBER	S1	S2	S 3	S4	S 5	S6	S 7	S8	S 9	S10
PARAMETER PCB 1016	77,000	ND	4,100	73,000	140,000	190,000	55,000	100,000	49,000	15,000
PCB 1254	45,000			170,000	i	<u>-</u>		550,000	83,000	
PCB 1260 TOTAL PCBs		2,000,000				-		730,000	···- ·· ·	61,000

ND = Non Detect

TABLE 4 ANALYTICAL DATA FOR DIOXINS AT THE SCRAP METAL SITE CHICAGO, ILLINOIS

Sample ID# Analyte	S1 μg/kg	S2 µg/kg	S3 µg/kg	S8 µg/kg	BLANK μg/kg
TOTAL 2378-TCDD EQUIVALENCE	0.265	4.004	1.207	0.36	0
2378-TCDF	0.22	ND	1.7*	ND	ND
TOTAL TCDF	1.5	17	9.4	1.9	ND
2378-TCDD	ND	ND	ND	ND	ND
TOTAL TCDD	0.034	0.58	0.13	0.15	ND
12378-PeCDF	0.094	4.4	0.55	ИD	ND
23478-PeCDF	0.26	6.3	1.3	0.22	ND
TOTAL PeCDF	1.3	42	7.6	3.3	ND
12378-PeCDD	ND	ND	ND	ND	ND
TOTAL PeCDD	ND	ND	ND	ND	ND
123478-HxCDF	0.25	ND	0.81	0.5	ND
123678-HxCDF	0.16	ND	0.64	ND	ND
234678-HxCDF	0.23	2.7	0.99	0.36	0.039
123789-HxCDF	0.095	0.890	0.22	0.17	ND
TOTAL HXCDF	2	15	7.7	2.7	0.039
123478-HxCDD	0.021*	ND	0.1	0.029	ND
123678-HxCDD	0.070*	0.25	0.17	0.670	ND
123789-HxCDD	0.033*	0.13	0.12	0.19	ND
TOTAL HXCDD	0.24	0.38	1.1	1.8	ND
1234678-HCDF	0.9	12	3	1.1	ND
1234789-HpCDF	0.12	3.6	0.33	0.24	ND.
TOTAL HPCDF	1.5	16	4.5	1.3	ND
1234678-HpCDD	0.73	2.9	1.5	2.8	0.0023
TOTAL HPCDD	1.4	5.9	2.9	5.7	0.0044
OCDF	0.77	30	1.6	1.6	ND
OCCD	4.3	22	3.9	51	0.021

ND - Non detect
* - Value may include contributions from other TCDF isomers.

ppb. The PCBs can be directly associated with past activities at the site as reported by the Heatbath employee and former railroad employee. On-site soil contain dioxins that were greater than 1 ug/kg (2,3,7,8-TCDD equivalence) S2 4.004 ug/kg and S3 1.207 ug/kg. CIEC workers continue work on-site where they have the potential to come in contact with hazardous substances.

6.2 Migration of Contaminants

On-site soils contains heavy metals that were above the RCRA limits for cadmium and lead (1 and 5 mg/L, respectively) which are listed wastes D006 and D008, respectively. Total PCBs were detected in on-site soils ranging from 37,100 to 2,000,000 ppb. On-site soils contain dioxins that were greater than 1 ug/kg (2,3,7,8-TCDD equivalence) in samples S2 4.004 ug/kg and S3 1.207 ug/kg. The Agency for Toxic Substances and Disease Registry (ATSDR) considers 1 ug/kg of dioxin in soil to be a level of concern in residential areas. The migration of contaminants from the facility is possible due to piles by airborne dusts or storm run-off which could effect on-site workers and nearby residential areas.

7.0 SUMMARY

The site assessment documented the existence of hazardous substances at the Scrap Metal site. These substances were detected in samples of on-site soils which were found to contain materials that possess characteristics of hazardous wastes, as defined by RCRA. The SM site is an active metal reclamation facility. Conditions at the SM site may pose risks to the business employees on-site and to nearby residential areas.

7.1 RCMS 4.0 COST PROJECTION

Removal cost estimates for off-site disposal were projected using the U.S. EPA Removal Cost Management System (RCMS) program. Cost estimates to mitigate threats due to on-site contaminated materials contained in the soil are based on the following assumptions:

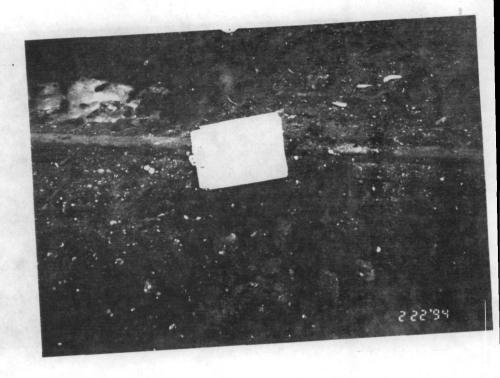
- 1. There is approximately 4,500 cubic yards contaminated soils on-site. This estimated amount was generated assuming that the 3-acre site could be contaminated to a depth of at least 1 feet. This estimate does not include the soils underneath the concrete in east lot.
- 2. Soil containing PCBs at concentrations grater than 50 ppm will be disposed of at a TSCA approved landfill;

- 3. Soil containing dioxin at concentrations grater than 1 ug/kg will be disposed of at a TSCA approved landfill;
- 4. Soil exhibiting TCLP cadmium and lead concentrations greater than 1 and 5 mg/l respectively will be stabilized either on-site or off-site and transported to an appropriate and RCRA permitted off-site landfill. If the soil also exhibits a PCB concentration greater than 50 mg/kg prior stabilization, the soil will be stabilized and disposed of at a TSCA approved landfill.

The cost estimates for the cleanup are based on 90 12-hour days by Emergency Response Cleanup Support (ERCS) contractor; Riedel Environmental Services, and is projected to cost \$1,696,050. See Appendix C for the RCMS cost estimate.

APPENDIX A PHOTO LOG

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SITE: SCRAP METAL .

DATE: 2-22-94

TIME: 1225

DIRECTION: NORTH

PHOTOGRAPHER: JOHN NORDINE

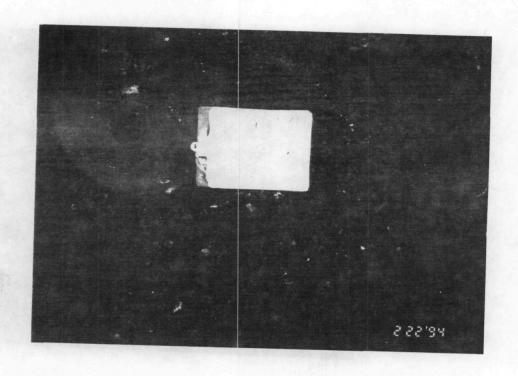
DESCR: S7 COLLECTED 25'N & 25'E OF THE

NW CORNER OF THE OFFICE



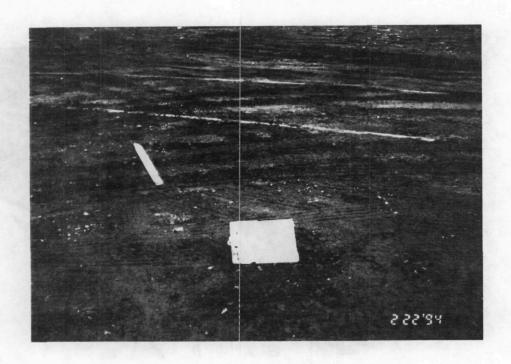
SITE: SCH.

DATE: 2-22
TIME: 1325
DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: S9 COLLECTED FROM THE MIDDLE OF
THE WEST LOT BY RAILROAD TRACKS SITE: SCRAP METAL DATE: 2-22-94 TIME: 1325

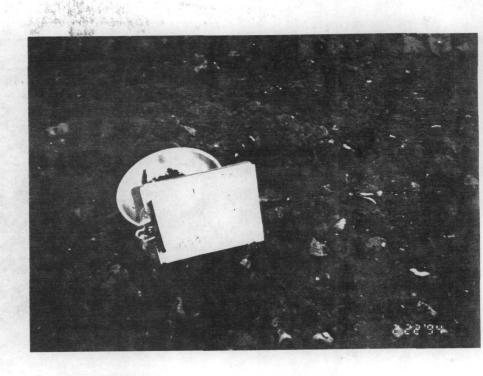


SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1325
DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE

DESCR: SP COLLECTED FROM THE MIDDLE OF THE WEST LOT BY RAILROAD TRACKS



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1250
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: S8 COLLECTED NEAR CRAIN ALONG
WEST PROPERTY LINE WEST LOT



SITE: SCRAP METAL .

DATE: 2-22-94

TIME: 1200

DIRECTION: NORTH

PHOTOGRAPHER: JOHN NORDINE

DESCR: S1 EAST LOT, 20' NORTH OF THE

OFFICES NE CORNER.

differently and the second



SITE: SCRAP METAL .

DATE: 2-22-94

TIME: 1203

DIRECTION: NORTH

PHOTOGRAPHER: JOHN NORDINE

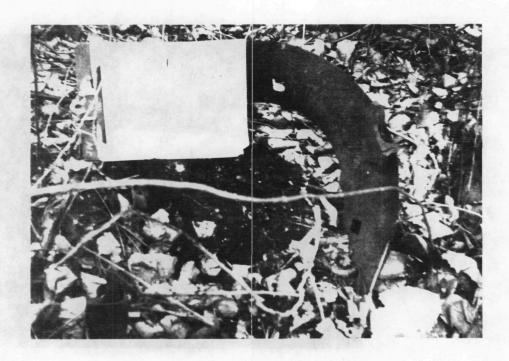
DESCR: S2 EAST LOT, WEST OF THE FORMER

WIRE INCINERATOR



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1203

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT, S2 LOCATED ALONG
RAILROAD OVERPASS



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1205
DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT SOIL SAMPLE S2
LOCATION



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1225

10 minds

DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: OIL STAINED AREA WERE S10 WAS
COLLECTED N SIDE OF THE OFFICE



SITE: SCRAP METAL .

DATE: 2-22-94
TIME: 1225
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: OIL STAINED AREA WERE S10 WAS
COLLECTED N SIDE OF THE OFFICE



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: NORTH SIDE OF THE OFFICE
BUILDING EAST LOT



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: NORTH SIDE OF THE OFFICE
BUILDING EAST LOT



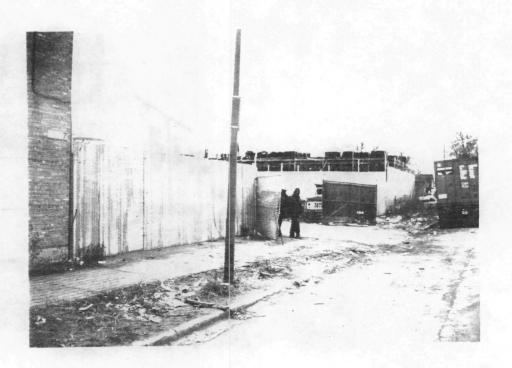
SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: MOBTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: FENCE ALONG THE WEST SIDE OF
THE EAST LOT

(秦)到1994年,《中国1994年



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: NORTH

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: NOTE FENCE ALONG WEST SIDE OF
THE OFFICE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1345 DIRECTION: SOUTH

PHOTOGRAPHER: JOHN NORDINE
DESCR: WEST SIDE OF THE OFFICE NOTE
FENCE



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1215 DIRECTION: NORTHWEST PHOTOGRAPHER: JOHN NORDINE DESCR: NORTHWEST AREA OF EAST LOT



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1230
DIRECTION: EAST

PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT CARDBOARD FIRE BY EAST
SIDE OF THE OFFICE BUILDING

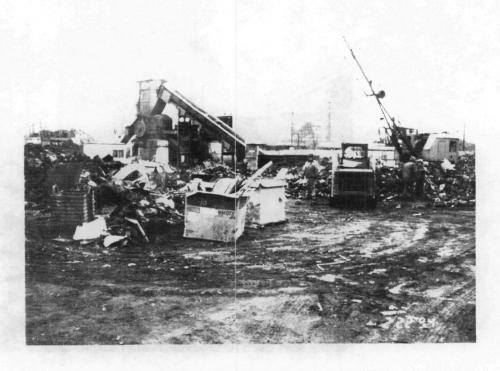


SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1210
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDI

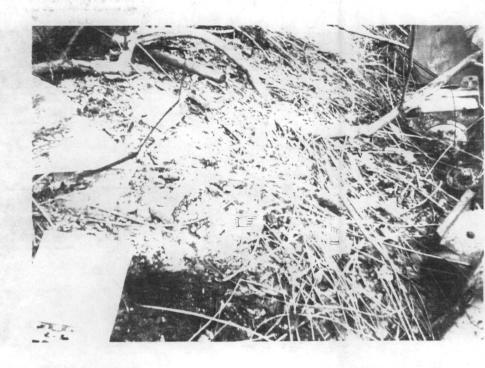
PHOTOGRAPHER: JOHN NORDINE
DESCR: NOTE SMOKE FROM CARDBOARD FIRE
LOCATED BY THE EAST LOT OFFICE



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST SIDE OF THE OFFICE
BUILDING EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1345 SITE: 50...
DATE: 2-22
TIME: 1345
DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: METAL SHREDDER NOTE DUST CLOUDS
EAST LOT



SITE: SCRAP METAL

DATE: 2-22-94

TIME: 1215

DIRECTION: NORTH

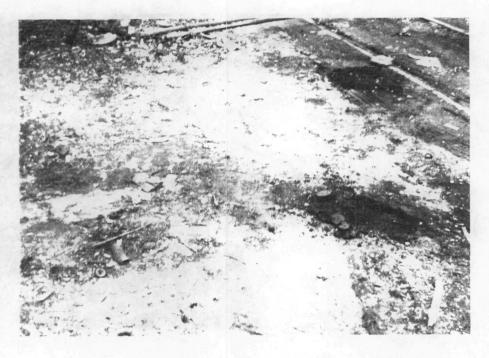
PHOTOGRAPHER: JOHN NORDINE

DESCR: SOIL SAMPLE S3 FROM ASH PILE IN

THE EAST LOT, NOTE THE DEAD RATS



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1305 DATE: DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: S4 COLLECTED 3' EAST OF THE
NORTHWEST GATE, WEST LOT



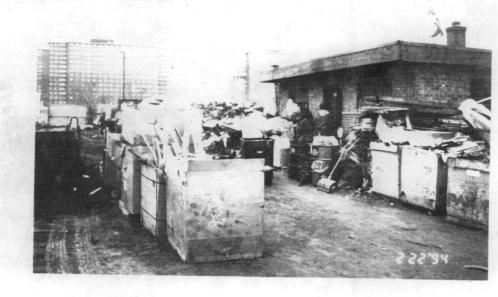
SITE: SCRAP METAL DATE: 2-22-94 TIME: 1305

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: S4 COLLECTED 3' EAST OF THE
NORTHWEST GATE, WEST LOT



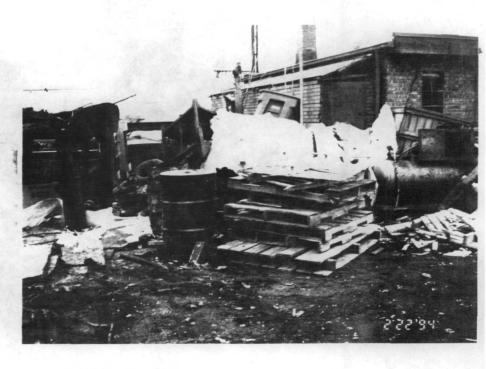
SITE: SCRAP METAL DATE: 2-22-94 TIME: 1310

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: S5 & S6 ASH PILE WEST LOT ALONG
NORTH PROPERTY LINE



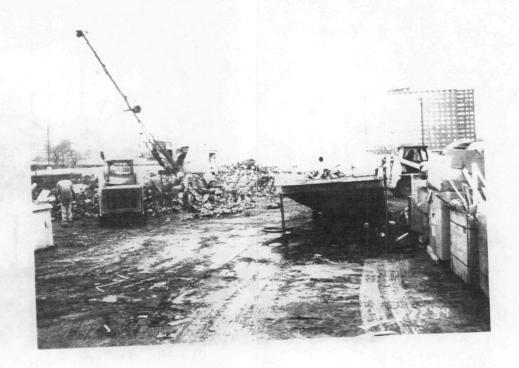
SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1230 DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT RECIEVING SHED FOR
RECYCLE MATERIALS

和"安徽"。



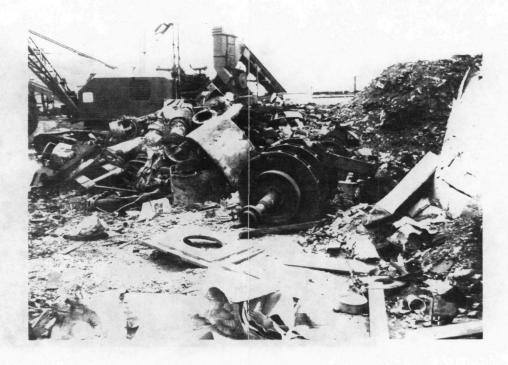
SITE: SCRAP METAL DATE: 2-22-94 TIME: 1230 DIRECTION: SOUTH

PHOTOGRAPHER: JOHN NORDINE DESCR: DRUM NEAR RECEIVING SHED CONTENTS KNOWN



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345

DIRECTION: EAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: SCRAP METAL PILES EAST LOT NOTE
APARTMENTS IN THE BACKGROUND



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1205 DIRECTION: EAST

PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT SCRAP METAL PILES



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1230

PA-AR

DIRECTION: WEST
PHOTOGRAPHER: JOHN NORDINE
DESCR: ALUMINUM FURNACES ALONG SOUTH
PROPERTY LINE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1345

DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: SCRAP METAL PILES AND ALUMINUM
FURNACES IN THE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1345 TIME:

DIRECTION: SOUTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: SOUTH PART OF THE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1345 DIRECTION: EAST PHOTOGRAPHER: JOHN NORDINE

DESCR: RECYCLING SHED IN THE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: , 1250 DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: NORTHEAST PART OF THE WEST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1250 DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: NORTHEAST PART OF THE WEST LOT



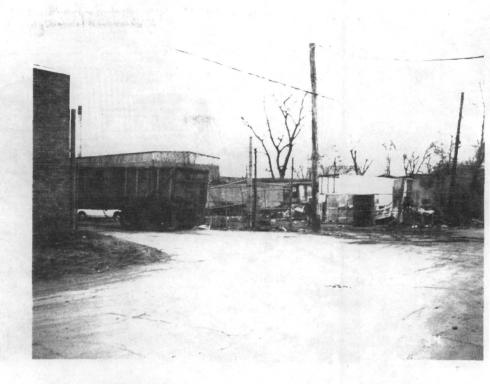
SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1250

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: CRAIN BOOM, ASH PILE, AND TIRES
ALONG NORTH SIDE OF WEST LOT

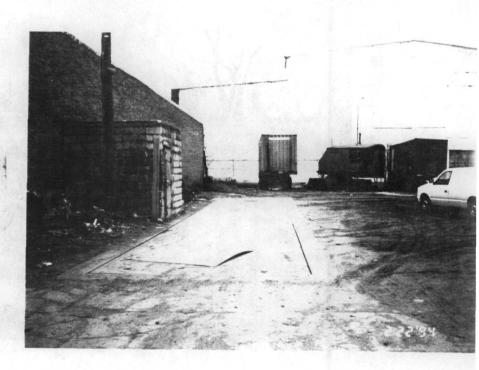


SITE: SCRAP METAL DATE: 2-22-94 TIME: 1305 DIRECTION: EAST

PHOTOGRAPHER: JOHN NORDINE
DESCR: BURN DRUM WEST LOT, NOTE BURNED
WIRE AT BASE OF THE DRUM



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1345 DIRECTION: NORTHWEST
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST SIDE OF THE WEST LOT, NOTE
FENCE AND GATES



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1325

DIRECTION: WEST
PHOTOGRAPHER: JOHN NORDINE
DESCR: SCALE SHED AND TRUCK SCALE IN
THE WEST LOT



SITE: SCRAP METAL .

DATE: 2-22-94
TIME: 1300
DIRECTION: SOUTHWEST
PHOTOGRAPHER: JOHN NORDINE
DESCR: SOUTHWEST PART OF THE WEST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1325 DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: NORTHSIDE OF WEST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1330 DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT SCRAP METAL PILES AND
EMPTY DRUMS



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345 DATE:

DIRECTION: EAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: FERROUS SORTER AND THE METAL
SHREDDER IN THE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1225 DIRECTION: WEST
PHOTOGRAPHER: JOHN NORDINE
DESCR: METAL SHREDDER AT THE EAST LOT



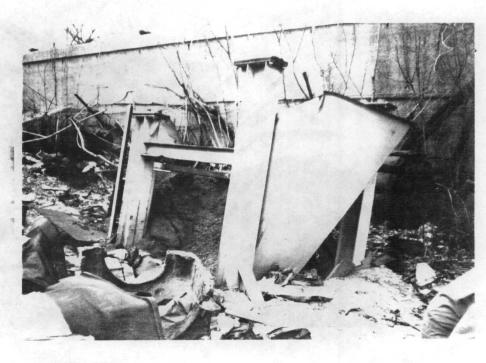
SITE: SCRAP METAL DATE: 2-22-94 TIME: 1330

DIRECTION: EAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: NOTE APARTMENTS IN BACK OF
SHREDDER & FERROUS SORTER

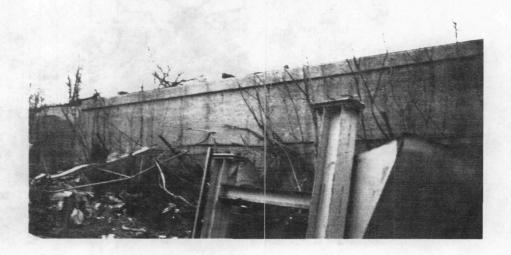


SITE: SCRAP METAL DATE: 2-22-94 TIME: 1201

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT RAILROAD OVERPASS
NORTH BOUNDARY OF SITE.



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1202 DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT PILES OF SCRAP METAL



SITE: SCRAP METAL

DATE: 2-22-94

TIME: 1202

DIRECTION: NORTH

PHOTOGRAPHER: JOHN NORDINE

DESCR: EAST LOT, PILES OF SCRAP METAL

BY RAILROAD OVERPASS



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1205

DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT SCRAP METAL PILES IN
THE NORTHEAST AREA

APPENDIX B DATA QUALITY ASSURANCE REVIEW





International Specialists in the Environment

111 West Jackson Boulevard Chicago, Illinois 60604

Tel: (312) 663-9415, Fax: (312) 663-0791

MEMORANDUM

DATE: May 3, 1994

TO: John Nordine, Project Manager, E&E, Chicago, IL

FROM: Nabil Fayoumi, TAT-Chemist, E&E, Chicago, IL NF

THRU: David Hendren, TAT-Chemist, E&E, Chicago, IL

SUBJ: Polychlorinated Biphenyl Data Quality Assurance Review, Scrap

Metal Site, Chicago, Cook County, Illinois.

REF: Analytical TDD: T059402807 , Project TDD: T059402007

Analytical PAN: EILO831AAA Project PAN: EILO831SAA

The data quality assurance review of 10 soil samples collected from the Scrap Metal Site in Chicago, Illinois has been completed. Analysis for Polychlorinated Biphenyls (PCBs) was performed by Twin City Testing located in St. Paul. Minnesota, Illinois in accordance with U.S. EPA Method SW-846-8080.

The samples were numbered as following:

TAT Sample #	Corresponding t	:0 -	->	Laboratory	Sample	#
S-1				15270		
s-2				15277		
s-3				15278		
s-4				15279		
S-5				15283		
s-6				15287		
s-7				15289		
S- 8				15290		
S-9				15291		
S-10				15292		

Data Qualifications:

I Sample Holding Time: Acceptable

The samples were collected on 2/22/94, extracted on 3/1/94, and analyzed on 3/6/94, and 3/7/94. The holding time criteria of 14 days for soil from collection to extraction was satisfied. The analysis of the samples was completed within the 40 day holding time requirement after extraction.

II Instrument Performance: Acceptable

The quality control criteria established for the surrogate retention time shift of less than 0.3% were acceptable. Peak resolution was adequate.

III Calibration: Qualified

A. Initial Calibration:

A 5-point calibration check was performed prior to sample analysis. The Relative Standard Deviations (RSDs) of all calibration factors for all Arochlors did not meet the quality control criteria of less than or equal to 10% RSD except 1260 on the DB-608 column, therefore; positive results were flagged (J) as estimated.

B. Continuing Calibration:

The percent difference (%D) requirements of less than 15% were acceptable for all data.

IV Method Blanks: Acceptable

A method blank was analyzed with the samples. No contaminants were detected above the instrument detection limit.

V Matrix Spike/Matrix Spike Duplicate: Acceptable

Spike Sample Analysis:

All Matrix Spike/Matrix Spike Duplicate recoveries were diluted out due to high level of PCBs in the samples. No action was necessary.

VI Compound Identification and Quantitation: Acceptable

All reported Retention Times were within the allowed RT windows. Positive results were confirmed using dual column analysis. The laboratory states in the narrative that because of the complexity of the Aroclor patterns present, it is possible that several Aroclors (PCB 1232 and PCB 1242) could be misidentified as PCB 1016.

VII Surrogate Recovery: Acceptable

The surrogates were diluted out due to high level of PCBs in the samples. No action was necessary.

VIII Overall Assessment of Data for Use

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality control Guidance For Removal Activities" (OSWER Directive 9360.4-01, April 1990).

Based upon the information provided, the data are acceptable for use with the above stated data qualifications.

Data Qualifiers and Definitions

J - The associated numerical value is an estimated quantity because the were less than the contract required detection limits or quality control criteria were not met.

• .

(All values are in μ g/Kg which is equal to parts-per-billion)

Client ID:

S-1

TCT ID:	15270	
Parameter:		POL
PCB 1016	77,000	7,800
PCB 1221	ND	7,800
PCB 1232	ND	7,800
PCB 1242	ND	7,800
PCB 1248	ND	7,800
PCB 1254	45,000	7,800
PCB 1260	21,000	7,800
% Surrogate #1 Recovery:	1 %	
% Surrogate #2 Recovery:	_1%	
Date Extracted:	3/1/94	

3/7/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Date Analyzed:

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:

S-2

TCT ID:	15277	
Parameter:		POL
PCB 1016	ND	250,000
PCB 1221	ND	250,000
PCB 1232	ND	250,000
PCB 1242	ND	250,000
PCB 1248	ND	250,000
PCB 1254	ND	250,000
PCB 1260	2,000,000	250,000
% Surrogate #1 Recovery:	1%	
% Surrogate #2 Recovery:	¹ %	
Date Extracted:	3/1/94	

3/7/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Date Analyzed:

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in μ g/Kg which is equal to parts-per-billion)

Client ID:	S-3
CHUIL ID.	55

TCT ID:	15278	
Parameter:		POL
PCB 1016	4,100	2,500
PCB 1221	ND	2,500
PCB 1232	ND	2,500
PCB 1242	ND	2,500
PCB 1248	ND	2,500
PCB 1254	18,000	2,500
PCB 1260	15,000	2,500
% Surrogate #1 Recovery:	1%	
% Surrogate #2 Recovery:	¹ %	
Date Extracted:	3/1/94	

¹Surrogate diluted out due to high level of PCB in sample.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Date Analyzed:

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference: EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.

3/7/94



(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:

S-4

TCT ID:	15279	
Parameter:		POL
PCB 1016	73,000	64,000
PCB 1221	ND	64,000
PCB 1232	ND	64,000
PCB 1242	ND	64,000
PCB 1248	ND	64,000
PCB 1254	160,000	64,000
PCB 1260	92,000	64,000
% Surrogate #1 Recovery:	¹ %	
% Surrogate #2 Recovery:	1 %	
Date Extracted:	3/1/94	

3/6/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Date Analyzed:

Surrogate #1 = TCMX (2,4,5,6-tetrachioro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition. Reference:



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in μ g/Kg which is equal to parts-per-billion)

Client ID:

S-5

TCT ID:	15283	
Parameter:		POL
PCB 1016	410,000	75,000
PCB 1221	ND	75,000
PCB 1232	ND	75,000
PCB 1242	ND	75,000
PCB 1248	ND	75,000
PCB 1254	510,000	75,000
PCB 1260	210,000	75,000
% Surrogate #1 Recovery:	¹ %	
% Surrogate #2 Recovery:	1%	
Data Entracted	2/14/04	

Date Extracted:

3/1/94

Date Analyzed:

3/6/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

CIL A ID	0.7
Client ID:	S-6

TCT ID:	15287

ICT ID:	13287	
Parameter:		POL
PCB 1016	190,000	72,000
PCB 1221	NĎ	72,000
PCB 1232	ND	72,000
PCB 1242	ND	72,000
PCB 1248	ND	72,000
PCB 1254	300,000	72,000
PCB 1260	200,000	72,000
% Surrogate #1 Recovery:	1 %	
% Surrogate #2 Recovery:	1 %	

Date Extracted: 3/1/94

3/6/94 Date Analyzed:

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference: EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client	m:
CHEIR	ш.

S-7

TCT ID:	15289	
Parameter:		<u>POL</u>
PCB 1016	55,000	13,000
PCB 1221	ND	13,000
PCB 1232	ND	13,000
PCB 1242	ND	13,000
PCB 1248	ND	13,000
PCB 1254	42,000	13,000
PCB 1260	39,000	13,000
% Surrogate #1 Recovery:	¹ %	
% Surrogate #2 Recovery:	¹ %	
Date Extracted:	3/1/94	
Date Analyzed:	3/7/94	

¹Surrogate diluted out due to high level of PCB in sample.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:

S-8

TCT ID:	15290	
Parameter:		POL
PCB 1016	$100,000^2$	75,000
PCB 1221	ND	75,000
PCB 1232	ND	75,000
PCB 1242	ND	75,000
PCB 1248	ND	75,000
PCB 1254	550,000	75,000
PCB 1260	80,000	75,000
% Surrogate #1 Recovery:	1%	
% Surrogate #2 Recovery:	1%	
Date Extracted:	3/1/94	
Date Analyzed:	3/7/94	

¹Surrogate diluted out due to high level of PCB in sample.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



²Value was not confirmed within 25%, the lower value is reported.

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:	S-9

TCT ID:

 Parameter:
 POL

 PCB 1016
 49,000
 14,000

PCB 1221 ND 14,000

15291

PCB 1232 ND 14,000

PCB 1242 ND 14,000

PCB 1248 ND 14,000

PCB 1254 83,000 14,000

PCB 1260 32,000 14,000

% Surrogate #1 Recovery: --1%

% Surrogate #2 Recovery: -1%

Date Extracted: 3/1/94

Date Analyzed: 3/7/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference: EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:

S-10

TCT ID:	15292	
Parameter:		<u>PQL</u>
PCB 1016	15,000	5,000
PCB 1221	ND	5,000
PCB 1232	ND	5,000
PCB 1242	ND	5,000
PCB 1248	ND	5,000
PCB 1254	24,000	5,000
PCB 1260	22,000	5,000
% Surrogate #1 Recovery:	1%	
% Surrogate #2 Recovery:	1%	

Date Extracted:

3/1/94

Date Analyzed:

3/7/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachioro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

ecology and environment, inc.



111 West Jackson Boulevard Chicago, Illinois 60604 Tel: (312) 663-9415, Fax: (312) 663-0791

MEMORANDUM

DATE: May 3, 1994

TO: John Nordine, Project Manager, E&E, Chicago, IL

FROM: Nabil Fayoumi, TAT-Chemist, E&E, Chicago, IL Nf

THRU: David Hendren, TAT-Chemist, E&E, Chicago, IL

SUBJ: Inorganic Data Quality Assurance Review, Scrap Metal Site,

Chicago, Cook County, Illinois.

REF: Analytical TDD: T059402807 Project TDD: T059402007

Analytical PAN: EIL0831AAA Project PAN: EIL0831SAA

The data quality assurance review of 10 Soil samples collected from the Scrap Metal Site in Chicago, Illinois has been completed. Analysis for Total and TCLP Metals was performed by Twin City Testing located in St. Paul, Minnesota in accordance with U.S. EPA Methods SW-846-6010/7000.

The samples were numbered as following:

TAT Sample #	Corresponding to ->	Laboratory Sample #
S-1	•	15270
S-2		15277
s-3		15278
S-4		15279
S-5		15283
s-6		15287
s-7		15289
s-8		15290
S -9		15291
s-10		15292

Data Qualifications:

I Sample Holding Time: Acceptable

The samples were collected on 2/22/94 and analyzed between 3/1/94, and 3/18/94. The holding time criteria of 6 months for metals and 28 days for mercury from collection to analysis was satisfied.

II Calibration: Acceptable

A. Initial Calibration:

Calibration results were within the established quality control limits of 90-110% of the true value for metals, and 80-120% for mercury.

B. Continuing Calibration:

Calibration results showed that the control criteria of 90-110% for metals, and 80-120% for mercury were satisfied.

III Method Blanks: Acceptable

A method blank was analyzed with the samples. No contaminants were detected above the instrument detection limit.

IV Interference Check Sample Analysis: Acceptable

All parameters were within the Interference Check Sample (ICS) control limits of 80-120% of the true values. ICS was run at the beginning and end of sample analysis.

V Matrix Spike/Matrix Spike Duplicate:

Spike Sample Analysis: Acceptable . .

All Matrix Spike/Matrix Spike Duplicate recoveries were within the control limits of 80-120%.

VI Optional Additional QC:

Laboratory Control Sample Analysis: Acceptable

The quality control criteria of 80-120% were met for the control sample.

VIII Overall Assessment of Data for Use

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality Control Guidance For Removal Activities" (OSWER Directive 9360.4-01, April 1990).

Based upon the information provided, the data are acceptable for use.

(All values are in μ g/L which is equivalent to parts-per-billion)

Client ID:

S1

S2

S3

TCT ID:

15270

15277

15278

<u>Parameter</u>				POL	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND	ND	100	3/15/94	6010
Barium	1,500	2,800	2,700	10	3/15/94	6010
Cadmium	870	1,300	110	10	3/15/94	6010
Chromium	ND	ND	ND	10	3/15/94	6010
Lead	7,500	16,000	690	50	3/15/94	6010
Mercury	ND	ND	ND	0.40	3/18/94	7470
Selenium	ND	ND	ND	100	3/15/94	6010
Silver	ND	ND	ND	10	3/15/94	6010
TCLP Date:	3/14/94	3/14/94	3/14/94			

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



(All values are in $\mu g/L$ which is equivalent to parts-per-billion, unless noted otherwise)

Client ID:

S4

S4

S4

TCT ID:

15279

15279D

15279S

<u>Parameter</u>				<u>PQL</u>	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND	106%	100	3/15/94	6010
Barium	1,100	1,100	80%	10	3/15/94	6010
Cadmium	230	220	90%	10	3/15/94	6010
Chromium	10	ND	91%	10	3/15/94	6010
Lead	8,200	8,100	81%	50	3/15/94	6010
Mercury	ND	ND	125%	0.40	3/18/94	7470
Selenium	ND	ND	113%	100	3/15/94	6010
Silver	ND	ND	92%	10	3/15/94	6010
TCLP Date:	3/14/94	3/14/94	3/14/94			

PQL = Practical Quantitation Limit

ND = Not Detected

D = Duplicate

S = Spike

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



(All values are in $\mu g/L$ which is equivalent to parts-per-billion)

Client ID:

S5

S6

S7

TCT ID:

15283

15287

15289

<u>Parameter</u>				POL	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND · -	ND	100	3/15/94	6010
Barium	4,800	2,700	1,300	10	3/15/94	6010
Cadmium	910	660	1,200	10	3/15/94	6010
Chromium	ND	ND	ND	10	3/15/94	6010
Lead	45,000	4,300	7,400	50	3/15/94	6010
Mercury	ND	ND	ND	0.40	3/18/94	7470
Selenium	ND	ND	ND	100	3/15/94	6010
Silver	ND	ND	ND	10	3/15/94	6010
TCLP Date:	3/14/94	3/14/94	3/14/94			

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



(All values are in µg/L which is equivalent to parts-per-billion)

Client ID: S8 S9 S10

TCT ID: 15290 15291 15292

<u>Parameter</u>				POL	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND · x	ND	100	3/15/94	6010
Barium	1,500	1,600	2,000	10	3/15/94	6010
Cadmium	850	480	690	10	3/15/94	6010
Chromium	12	15	ND	10	3/15/94	6010
Lead	5,300	71,000	32,000	50	3/15/94	6010
Mercury	ND	ND	ND	0.40	3/18/94	7470
Selenium	ND	ND	ND	100	3/15/94	6010
Silver	ND	ND	ND	10	3/15/94	6010
TCLP Date:	3/14/94	3/14/94	3/14/94			

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.





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International Specialists in the Environment

MEMORANDUM

DATE: March 29, 1994

TO: John Nordine, Project Manager, E & E, Chicago, IL

FROM: Nabil Fayoumi, TAT-Chemist, E & E, Chicago, IL

THRU: David Hendren, TAT-Chemist, E & E, Chicago, IL

SUBJ: Organic Data Quality Assurance Review, Scrap Metal Site,

Chicago, Cook County, Illinois.

REF: Analytical TDD: T059402807 '~ Project TDD: T059402007

Analytical PAN: EILO831AAA Project PAN: EILO831SAA

The data quality assurance review of 10 soil samples collected from the Scrap Metal Site in Chicago, Illinois has been completed. Analysis for Semivolatile Organics (SVOA) was performed by Twin City Testing located in St. Paul, Minnesota in accordance with U.S. EPA SW-846 method 8270.

The samples were numbered as following:

TAT Sample 🗲	Corresponding to ->	Laboratory Sample	£
S-1		15270	
S-2		15277	
s-3		15278	
S-4		15279	
S-5		15283	
S-6		15287	
S-7		15289	
5- 8		15290	
8-9		15291	
5-10		15292	

Data Qualifications:

I Holding Time: Acceptable

The samples were collected on 2/22/94, extracted on 3/3/94, and analyzed between 3/8/94 and 3/10/94. The holding time criteria from collection to extraction of 14 days was satisfied. Extracts were analyzed within 40 days.

II GC/MS Tuning: Acceptable

GC/MS ion abundance criteria using Decafluorotriphenylphosphine (DFTPP) were acceptable.

III Calibration: Acceptable

A. Initial Calibration:

A 5-point initial calibration was performed prior to analysis. All average relative response factors were greater than 0.05. All percent relative standard deviation (%RSD) between response factors were less than 30%, except 2,4-dinitrophenol (33.1 % RSD), which was not detected in any sample; therefore qualification was not necessary.

B. Continuing Calibration:

The percent difference (%D) between initial and continuing calibration were within the quality control criteria of less than or equal to 25% for all detected compounds.

IV Method Blank: Acceptable

The method blank was contaminated with Di-n-butylphthalate (47 PPB). No action was necessary because all reported sample values were greater than 10 times blank concentration level.

V Surrogate Recovery: Acceptable

Surrogates in the matrix spike sample (number S-4MS) were recovered at less than 10% recovery. Since both samples S-4 and S-4MSD displayed acceptable surrogate recoveries, it appears that the poor surrogate response in sample S-4MS resulted from sample preparation problems.

VI Matrix Spike/Matrix Spike Duplicates: Not Applicable

Fourteen out of 22 percent recoveries and 10 out of 11 relative percent differences for the MS/MSD were outside the established quality control criteria. This again probably resulted from sample S-4MS preparation problems.

VII Internal Standards: Qualified

The internal standard (chrysene D-12) area counts were outside (below) the range of -50% to +100% from the associated calibration standard for the following samples S-3, S-4, S-5, S-6, S-7, S-8, S-9, and S-10; therefore, all positive values quantitated from chrysene D-12 in these samples are flagged "J" (estimated).

VIII Overall Assessment of Data for Use

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality Control Guidance for Removal Activities" (OSWER 9360.4-01 April, 1990). Based upon the information provided, the data are acceptable for use with the above stated qualifications.

Data Qualifiers and Definitions:

J - The associated numerical value is an estimated quantity because the reported concentration is less than the contract required detection limit or quality control criteria were not met.

TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Client ID: S1 Lab ID (HSN): 15270
Matrix: SOIL Filename: 4068K08

Date Sampled: 02/22/94 Sample Size: 15.5 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL

Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 24.08

Compounds:	ug/Kg (PPB)	EQL
Phenol	1700	1700
bis(2-Chloroethyl)ether	1700 U	1700
2-Chlorophenol	1700 U	1700
1,3-Dichlorobenzene	1700 U	1700
1,4-Dichlorobenzene	1700 U	1700
1,2-Dichlorobenzene	1700 U	1700
2-Methylphenol	1700 ប	1700
2,2'-oxybis(1-Chloropropane)	1700 U	1700
4-Methylphenol	220 J	1700
N-Nitroso-di-n-propylamine	1700 U	1700
Hexachloroethane	1700 U	1700
Nitrobenzene	1700 U	1700
Isophorone	1700 U	1700
2-Nitrophenol	1700 U	1700
2,4-Dimethylphenol	1700 U	1700
bis(2-Chloroethoxy)methane	1700 U	1700
2,4-Dichlorophenol	1700 U	1700
1,2,4-Trichlorobenzene	1300 J	1700
Naphthalene	9 7 0 J	1700
4-Chloroaniline	1700 U	1700
Hexachlorobutadiene	1700 U	1700
4-Chloro-3-methylphenol	1700 U	1700
2-Methylnaphthalene	2900	1700
Hexachlorocyclopentadiene	1700 U	1700
2,4,6-Trichlorophenol	1700 U	1700
2,4,5-Trichlorophenol	4200 U	4200
2-Chloronaphthalene	1700 U	1700
2-Nitroaniline	4200 U	4200
Dimethylphthalate	510 J	1700
Acenaphthylene	1700 U	1700
2,6-Dinitrotoluene	1700 U	1700
3-Nitro aniline Acenaphth ene	4200 U 710 J	4200 1700
2,4-Dinitrophenol		4200
	4200 U	4200
4-Nitrophenol Dibenzofuran	1700 U	1700
2,4-Dinitrotoluene	1700 U	1700
Diethylphthalate	1700 U	1700
4-Chlorophenyl-phenylether	1700 U	1700
Fluorene	1700 J	1700
4-Nitroaniline	4200 U	4200
4,6-Dinitro-2-methylphenol	4200 U	4200
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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS (CONTINUED) EPA METHOD 8270

Client ID: S1
Matrix: SOIL

Lab ID (HSN): 15270 Filename: 4068K08

Compounds:	ug/Kg (PI	PB)	EQL
N-Nitrosodiphenylamine	1700	U	1700
4-Bromophenyl-phenylether	1700	U	1700
Hexachlorobenzene	1700	IJ	1700
Pentachlorophenol	4200	U	4200
Phenanthrene	3900		1700
Anthracene	290	J	1700
Carbazole	1700	U	1700
Di-n-butylphthalate	2600	В	1700
Fluoranthene	2500		1700
Pyrene	4200		1700
Butylbenzylphthalate	2500		1700
3,3'-Dichlorobenzidine	1700	U	1700
Benz(a)anthracene	1700	J	1700
Chrysene	1900		1700
bis(2-Ethylhexyl)phthalate	100000	E	1700
Di-n-octylphthalate	1700	U	1700
Benzo(b)fluoranthene	2800		1700
Benzo(k) fluoranthene	1100	J	1700
Benzo(a)pyrene	2000		1700
Indeno(1,2,3-cd)pyrene	1600	J	1700
Dibenz(a,h)anthracene	1700	U	1700
Benzo(g,h,i)perylene	1900		1700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	55%	25-121%
Phenol-d5	63%	24-113%
2-Chlorophenol-d4	76%	20-130%
1,2-Dichlorobenzene-d4	87%	20-130%
Nitrobenzene-d5	51%	23-120%
2-Fluorobiphenyl	114%	30-115%
2,4,6-Tribromophenol	57%	19-122%
Terphenyl-d14	112%	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Compounds:	ug/Kg (PPB)	EQL
Phenol	3100 JD	17000
bis(2-Chloroethyl)ether	17000 UD	17000
2-Chlorophenol	17000 UD	17000
1,3-Dichlorobenzene	17000 UD	17000
1,4-Dichlorobenzene	17000 UD	17000
1,2-Dichlorobenzene	17000 UD	17000
2-Methylphenol	17000 UD	17000
2,2'-oxybis(1-Chloropropane)	17000 UD	17000
4-Methylphenol	17000 UD	17000
N-Nitroso-di-n-propylamine	17000 UD	17000
Hexachloroethane	17000 UD	17000
Nitrobenzene '7	17000 UD	17000
Isophorone	17000 UD	17000
2-Nitrophenol	17000 UD	17000
2,4-Dimethylphenol	17000 UD	17000
bis(2-Chloroethoxy)methane	17000 UD	17000
2,4-Dichlorophenol	17000 UD	17000
1,2,4-Trichlorobenzene	17000 UD	17000
Naphthalene	17000 UD	17000
4-Chloroaniline	17000 UD	17000
Hexachlorobutadiene	17000 UD	17000
4-Chloro-3-methylphenol	17000 UD	17000
2-Methylnaphthalene	3200 JD	17000
Hexachlorocyclopentadiene	17000 UD	17000
2,4,6-Trichlorophenol	17000 UD	17000
2,4,5-Trichlorophenol	42000 UD	42000
2-Chloronaphthalene	17000 UD	17000
2-Nitroaniline	42000 UD	42000
Dimethylphthalate	17000 UD	17000
Acenaphthylene	17000 UD	17000
2,6-Dinitrotoluene	17000 UD	17000
3-Nitro aniline	42000 UD	42000
Acenaphth ene	17000 UD	17000
2,4-Dinitrophenol	42000 UD	42000
4-Nitrophenol	42000 UD	42000
Dibenzofuran	17000 UD	17000
2,4-Dinitrotoluene	17000 UD	17000
Diethylphthalate	17000 UD	17000
4-Chlorophenyl-phenylether	17000 UD	17000
Fluorene	1700 JD	17000
4-Nitroaniline	42000 UD	42000
4,6-Dinitro-2-methylphenol	42000 UD	42000
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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS (CONTINUED) EPA METHOD 8270

Client ID: S1 Lab ID (HSN): 15270 DL Matrix: SOIL Filename: 4067L15

Compounds:	ug/Kg (PF	PB) EQL
N-Nitrosodiphenylamine	17000	UD 17000
4-Bromophenyl-phenylether	17000	UD 17000
Hexachlorobenzene	17000	UD 17000
Pentachlorophenol	42000	UD 42000
Phenanthrene	4100	JD 17000
Anthracene	17000	UD 17000
Carbazole	17000	UD 17000
Di-n-butylphthalate	4700	JDB 17000
Fluoranthene	2900	JD 17000
Pyrene	7100	JD 17000
Butylbenzylphthalate	7700	JD 17000
3,3'-Dichlorobenzidine	17000	UD 17000
Benz(a)anthracene	1800	JD 17000
Chrysene	2000	JD 17000
bis(2-Ethylhexyl)phthalate	200000	ED 17000
Di-n-octylphthalate	17000	UD 17000
Benzo(b) fluoranthene	3000	JD 17000
Benzo(k)fluoranthene	17000	UD 17000
Benzo(a)pyrene	1900	JD 17000
Indeno(1,2,3-cd)pyrene	17000	UD 17000
Dibenz(a,h)anthracene	17000	UD 17000
Benzo(g,h,i)perylene	17000	UD 17000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	79 % JD	25-121%
Phenol-d5	93 %J D	24-113%
2-Chlorophenol-d4	89 %J D	20-130%
1,2-Dichlorobenzene-d4	90 % JD	20-130%
Nitrobenzene-d5	90 %J D	23-120%
2-Fluorobiphenyl	116 % JD	30-115%
2,4,6-Tribromophenol	73%J D	19-122%
Terphenyl-d14	152%JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

= Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Client ID: S1 Lab ID (HSN): 15270 DL2
Matrix: SOIL Filename: 4068K04

Date Sampled: 02/22/94 Sample Size: 15.5 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL
Date Extracted: 03/03/94 Dil. Factor: 20
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 24.08

Phenol 34000 UD 34000 2-Chlorophenol 34000 UD 34000 34000 UD 3400	Compounds:	ug/Kg (PPB)	EQL
2,4,6-Trichlorophenol 34000 UD 34000 2,4,5-Trichlorophenol 85000 UD 85000 2-Chloronaphthalene 34000 UD 34000 2-Nitroaniline 85000 UD 34000 Dimethylphthalate 34000 UD 34000 Acenaphthylene 34000 UD 34000 2,6-Dinitrotoluene 34000 UD 34000 3-Nitroaniline 85000 UD 85000 Acenaphthene 34000 UD 34000 2,4-Dinitrophenol 85000 UD 85000 4-Nitrophenol 85000 UD 85000 Dibenzofuran 34000 UD 34000 2,4-Dinitrotoluene 34000 UD 34000 Diethylphthalate 34000 UD 34000 4-Chlorophenyl-phenylether 34000 UD 34000	Phenol bis (2-Chloroethyl) ether 2-Chlorophenol 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene 2-Methylphenol 2,2'-oxybis (1-Chloropropane) 4-Methylphenol N-Nitroso-di-n-propylamine Hexachloroethane Nitrobenzene Isophorone 2-Nitrophenol 2,4-Dimethylphenol bis (2-Chloroethoxy) methane 2,4-Dichlorophenol 1,2,4-Trichlorobenzene Naphthalene 4-Chloroaniline Hexachlorobutadiene 4-Chloro-3-methylphenol 2-Methylnaphthalene	34000 UD	34000 34000
4-Nitroaniline 85000 UD 85000 4,6-Dinitro-2-methylphenol 85000 UD 85000	2-Methylnaphthalene Hexachlorocyclopentadiene 2,4,6-Trichlorophenol 2,4,5-Trichlorophenol 2-Chloronaphthalene 2-Nitroaniline Dimethylphthalate Acenaphthylene 2,6-Dinitrotoluene 3-Nitroaniline Acenaphthene 2,4-Dinitrophenol 4-Nitrophenol Dibenzofuran 2,4-Dinitrotoluene Diethylphthalate 4-Chlorophenyl-phenylether Fluorene 4-Nitroaniline	34000 UD 34000 UD 34000 UD 85000 UD 34000 UD	34000 34000 85000 34000 85000 34000 34000 85000 85000 85000 34000 34000 34000 34000

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Client ID: S1 Lab ID (HSN): 15270 DL2
Matrix: SOIL Filename: 4068K04

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	34000 UD	34000
4-Bromophenyl-phenylether	34000 UD	34000
Hexachlorobenzene	34000 UD	34000
Pentachlorophenol	85000 UD	85000
Phenanthrene	4600 JD	34000
Anthracene	34000 UD	34000
Carbazole	34000 UD	34000
Di-n-butylphthalate	4900 JDB	34000
Fluoranthene	3700 JD	34000
Pyrene	8000 JD	34000
Butylbenzylphthalate	8500 JD	34000
3,3'-Dichlorobenzidine	34000 UD	34000
Benz(a)anthracene	34000 UD	34000
Chrysene	34000 UD	34000
bis(2-Ethylhexyl)phthalate	180000 D	34000
Di-n-octylphthalate	34000 UD	34000
Benzo(b) fluoranthene	34000 UD	34000
Benzo(k)fluoranthene	34000 UD	34000
Benzo(a)pyrene	34000 UD	34000
Indeno(1,2,3-cd)pyrene	34000 UD	34000
Dibenz(a,h)anthracene	34000 UD	34000
Benzo(g,h,i)perylene	34000 UD	34000

Surrogate Recovery 2-Fluorophenol Phenol-d5 2-Chlorophenol-d4 1,2-Dichlorobenzene-d4 Nitrobenzene-d5 2-Fluorobiphenyl	85%JD 101%JD 91%JD 92%JD 108%JD 112%JD	QC LIMITS 25-121% 24-113% 20-130% 20-130% 23-120% 30-115%
2-Fluorobiphenyl 2,4,6-Tribromophenol Terphenyl-d14		

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S2 Lab ID (HSN): 15277 Matrix: SOIL Filename: 4068K09

15.2 grams Date Sampled: 02/22/94 Sample Size: Extract Vol.:
Dil. Factor:
GPC Factor: Date Received: 02/24/94 1000 uL Date Extracted: 03/03/94 1 Date Analyzed: 03/09/94

2 % Moisture: 28.24

Compounds:	ug/Kg (PPB)	EQL
Phenol	1800 U	1800
bis(2-Chloroethyl)ether	1800 U	1800
2-Chlorophenol	1800 U	1800
1,3-Dichlorobenzene	1800 U	1800
1,4-Dichlorobenzene	1800 U	1800
1,2-Dichlorobenzene	1800 U	1800
2-Methylphenol	1800 U	1800
2,2'-oxybis(1-Chloropropane)	1800 U	1800
4-Methylphenol	1800 U	1800
N-Nitroso-di-n-propylamine	1800 U	1800
Hexachloroethane	1800 U	1800
Nitrobenzene	1800 U	1800
Isophorone	1800 U	1800
2-Nitrophenol	1800 U	1800
2,4-Dimethylphenol	1800 U	1800
bis(2-Chloroethoxy)methane	1800 U	1800
2,4-Dichlorophenol	1800 U	1800
1,2,4-Trichlorobenzene	7500	1800
Naphthalene	180 J	1800
4-Chloroaniline	1800 U	1800
Hexachlorobutadiene	1800 U	1800
4-Chloro-3-methylphenol	1800 U	1800
2-Methylnaphthalene	180 J	1800
Hexachlorocyclopentadiene	1800 U	1800
2,4,6-Trichlorophenol	1800 U	1800
2,4,5-Trichlorophenol	4600 U	4600
2-Chloronaphthalene	1800 U	1800
2-Nitroaniline	4600 U	4600
Dimethylphthalate	1800 U	1800
Acenaphthylene	1800 U	1800
2,6-Dinitrotoluene	1800 U	1800
3-Nitroaniline	4600 U	4600
Acenaphthene	1800 U	1800
2,4-Dinitrophenol	4600 U	4600
4-Nitrophenol	4600 U	4600
Dibenzofuran	1800 U	1800
2,4-Dinitrotoluene	1800 U	1800
Diethylphthalate	1800 U	1800
4-Chlorophenyl-phenylether	1800 U	1800
Fluorene	1800 U	1800
4-Nitroaniline	4600 U	4600
4,6-Dinitro-2-methylphenol	4600 U	4600

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Client ID: S2
Matrix: SOIL

Lab ID (HSN): 15277 Filename: 4068K09

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	1800 U	1800
4-Bromophenyl-phenylether	1800 U	1800
Hexachlorobenzene	1800 U	1800
Pentachlorophenol	4600 U	4600
Phenanthrene	1300 J	1800
Anthracene	180 J	1800
Carbazole	1800 U	1800
Di-n-butylphthalate	2200 B	1800
Fluoranthene	1200 J	1800
Pyrene	2100	1800
Butylbenzylphthalate	1800 U	1800
3,3'-Dichlorobenzidine	1800 U	1800
Benz(a)anthracene	1100 J	1800
Chrysene	1400 J	1800
bis(2-Ethylhexyl)phthalate	1200 J	1800
Di-n-octylphthalate	1800 U	1800
Benzo(b) fluoranthene	1600 J	1800
Benzo(k)fluoranthene	940 J	1800
Benzo(a)pyrene	1200 J	1800
Indeno(1,2,3-cd)pyrene	1500 J	1800
Dibenz(a,h)anthracene	1800 U	1800
Benzo(g,h,i)perylene	1700 J	1800

Surrogate Recovery		QC LIMITS
2-Fluorophenol	36 %	25-121%
Phenol-d5	38%	24-113%
2-Chlorophenol-d4	478	20-130%
1,2-Dichlorobenzene-d4	56 %	20-130%
Nitrobenzene-d5	33 % J	23-120%
2-Fluorobiphenyl	61%	30-115%
2,4,6-Tribromophenol	32%	19-122%
Terphenyl-d14	70%	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S2 Lab ID (HSN): 15277 DL Matrix: SOIL Filename: 4067L16

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL

Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 28.24

% Moisture: 28.24

Compounds:	ug/Kg (PPB)	EQL
Phenol bis(2-Chloroethyl)ether 2-Chlorophenol 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene 2-Methylphenol 2,2'-oxybis(1-Chloropropane) 4-Methylphenol N-Nitroso-di-n-propylamine Hexachloroethane Nitrobenzene Isophorone 2-Nitrophenol 2,4-Dimethylphenol	18000 18000 18000 18000 18000 18000 18000 18000 18000 18000 18000		18000 18000 18000 18000 18000 18000 18000 18000 18000 18000 18000 18000
<pre>bis(2-Chloroethoxy)methane 2,4-Dichlorophenol 1,2,4-Trichlorobenzene</pre>		0 JD 0 UD	18000 18000 18000
Naphthalene 4-Chloroaniline Hexachlorobutadiene 4-Chloro-3-methylphenol	18000 18000 18000	0 UD	18000 18000 18000 18000
2-Methylnaphthalene Hexachlorocyclopentadiene 2,4,6-Trichlorophenol	18000 18000 18000	0 UD 0 UD 0 UD	18000 18000 18000
2,4,5-Trichlorophenol 2-Chloronaphthalene 2-Nitroaniline Dimethylphthalate Acenaphthylene 2,6-Dinitrotoluene	4600 1800 4600 1800 1800	0 UD 0 UD 0 UD 0 UD	46000 18000 46000 . 18000 18000
3-Nitroaniline Acenaphth ene 2,4-Dinitrophenol 4-Nitrophenol	46000 18000 46000	0 UD 0 UD 0 UD 0 UD	46000 18000 46000 46000
Dibenzofuran 2,4-Dinitrotoluene Diethylphthalate 4-Chlorophenyl-phenylether Fluorene 4-Nitroaniline 4,6-Dinitro-2-methylphenol	1800 1800 1800 1800 1800 4600	0 UD 0 UD 0 UD 0 UD 0 UD	18000 18000 18000 18000 18000 46000
1,0 Dinicio-2-meenyiphenoi	-2000		40000

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Client ID: S2
Matrix: SOIL

Lab ID (HSN): 15277 DL Filename: 4067L16

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine 4-Bromophenyl-phenylether Hexachlorobenzene Pentachlorophenol Phenanthrene Anthracene Carbazole Di-n-butylphthalate Fluoranthene Pyrene Butylbenzylphthalate 3,3'-Dichlorobenzidine Benz(a)anthracene Chrysene bis(2-Ethylhexyl)phthalate Di-n-octylphthalate	18000 UD 3000 JDB 18000 UD 2800 JD 18000 UD	18000 18000 18000 18000 18000 18000 18000 18000 18000 18000 18000 18000 18000
Benzo(b) fluoranthene Benzo(k) fluoranthene Benzo(a) pyrene Indeno(1,2,3-cd) pyrene Dibenz(a,h) anthracene Benzo(g,h,i) perylene	2100 JD 18000 UD 18000 UD 18000 UD 18000 UD 18000 UD	18000 18000 18000 18000 18000

Surrogate Recovery 2-Fluorophenol Phenol-d5 2-Chlorophenol-d4 1,2-Dichlorobenzene-d4 Nitrobenzene-d5 2-Fluorobiphenyl 2,4,6-Tribromophenol	49%JD 56%JD 53%JD 56%JD 56%JD 66%JD 36%JD	QC LIMITS 25-121% 24-113% 20-130% 20-130% 23-120% 30-115% 19-122%
Terphenyl-d14	90 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S3 Lab ID (HSN): 15278
Matrix: SOIL Filename: 4068K10

Date Sampled: 02/22/94 Sample Size: 15 grams
Date Received: 02/24/94 Extract Vol.: 500 uL
Date Extracted: 03/03/94 Dil. Factor: 1

Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 30.97

Compounds:	ug/Kg (PPB)	EQL
Phenol	97	ט 0'	970
bis(2-Chloroethyl)ether		'0 U	970
2-Chlorophenol		'0 U	970
1,3-Dichlorobenzene		'0 U	970
1,4-Dichlorobenzene		'0 U	970
1,2-Dichlorobenzene		'0 Ū	970
2-Methylphenol		'0 U	970
2,2'-oxybis(1-Chloropropane)		'0 U	970
4-Methylphenol		'0 U	970
N-Nitroso-di-n-propylamine		'0 Ü	970
Hexachloroethane		'0 Ü	970
Nitrobenzene		'0 U	970
Isophorone		'0 U	970
2-Nitrophenol		'O U	970
2,4-Dimethylphenol		'0 Ŭ	970
bis (2-Chloroethoxy) methane		'O Ü	970
2,4-Dichlorophenol		'0 Ü	970
1,2,4-Trichlorobenzene		'0 Ü	970
Naphthalene		'0 U	970
4-Chloroaniline		'0 U	970
Hexachlorobutadiene		'0 U	970
4-Chloro-3-methylphenol		ט 0'	970
2-Methylnaphthalene		'0 U	970
Hexachlorocyclopentadiene		'0 U	970
2,4,6-Trichlorophenol		'0 Ū	970
2,4,5-Trichlorophenol			2400
2-Chloronaphthalene		'0 U	970
2-Nitroaniline			2400
Dimethylphthalate		'0 U	970
Acenaphthylene	97	'0 U	970
2,6-Dinitrotoluene	97	'O U	970
3-Nitroaniline	240	U 0	2400
Acenaphth ene	97	'O U	970
2,4-Dinitrophenol	³ 240	0 U	2400
4-Nitrophenol	240	U 0	2400
Dibenzofuran	97	'O U	970
2,4-Dinitrotoluene	97	'O U	970
Diethylphthalate	97	'O U	970
4-Chlorophenyl-phenylether	97	'O U	970
Fluorene	97	'O U	970
4-Nitroaniline	240	U 0	2400
4,6-Dinitro-2-methylphenol	240	0 U	2400
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Client ID: S3
Matrix: SOIL

Lab ID (HSN): 15278 Filename: 4068K10

Compounds:	ug/Kg (PPB)	EQL
-		
N-Nitrosodiphenylamine	9 7 0 U	970
4-Bromophenyl-phenylether	9 7 0 U	970
Hexachlorobenzene	970 U	970
Pentachlorophenol	2400 U	2400
Phenanthrene	880 J	970
Anthracene	140 J	970
Carbazole	970 Ŭ	9 7 0
Di-n-butylphthalate	1100 B	970
Fluoranthene	1700	970
Pyrene	3500 ケ Y	970
Butylbenzylphthalate	970 U Y	970
3,3'-Dichlorobenzidine	970 U Y	970
Benz(a)anthracene	2000 T Y	970
Chrysene	2200 T Y	970
bis(2-Ethylhexyl)phthalate	3300 ブ Y	970
Di-n-octylphthalate	970 U	970
Benzo(b) fluoranthene	3600	970
Benzo(k) fluoranthene	1100	970
Benzo(a) pyrene	2600	970
Indeno(1,2,3-cd)pyrene	2600	970
Dibenz(a,h)anthracene	970 U	970
Benzo(g,h,i)perylene	2600	970

Surrogate Recovery		QC LIMITS
2-Fluorophenol	25%	25-121%
Phenol-d5	34%	24-113%
2-Chlorophenol-d4	38%	20-130%
1,2-Dichlorobenzene-d4	40%	20-130%
Nitrobenzene-d5	29%	23-120%
2-Fluorobiphenyl	64%	30-115%
2,4,6-Tribromophenol	41%	19-122%
Terphenyl-d14	84%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

f = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S3 Lab ID (HSN): 15278 DL Matrix: SOIL Filename: 4067L17

Date Sampled: 02/22/94 Sample Size: 15 grams
Date Received: 02/24/94 Extract Vol.: 500 uL
Date Extracted: 03/03/94 Dil. Factor: 10

Date Analyzed: 03/09/94 GPC Factor: 2 % Moisture: 30.97

Compounds:	ug/Kg (1	PPB)	EQL
Phenol	9700	מט כ	9700
bis(2-Chloroethyl)ether		סט ס	9700
2-Chlorophenol		סס ס	9700
1,3-Dichlorobenzene		טס (9700
1,4-Dichlorobenzene		סט כ	9700
1,2-Dichlorobenzene		סט ס	9700
2-Methylphenol		סט ס	9700
2,2'-oxybis(1-Chloropropane)) UD	9700
4-Methylphenol		סט כ	9700
N-Nitroso-di-n-propylamine	9700) UD	9700
Hexachloroethane		סט כ	9700
Nitrobenzene		טס (9700
Isophorone		ָ סס	9700
2-Nitrophenol		סט ס	9700
2,4-Dimethylphenol		סט ס	
bis(2-Chloroethoxy)methane		סס ס	
2,4-Dichlorophenol		סט כ	
1,2,4-Trichlorobenzene		סט ס	
Naphthalene		סט ס	
4-Chloroaniline		טס ס	
Hexachlorobutadiene		טס ס	9700
4-Chloro-3-methylphenol		ט ס	9700
2-Methylnaphthalene		סט ס	9700
Hexachlorocyclopentadiene		סט ס	9700
2,4,6-Trichlorophenol		מט מ	9700
2,4,5-Trichlorophenol	24000		24000
2-Chloronaphthalene		סס ס	9700
2-Nitroaniline	24000		24000
Dimethylphthalate		סט ס	9700
Acenaphthylene		סט ס	9700
2,6-Dinitrotoluene		OD (9700
3-Nitroaniline	24000	UD (24000
Acenaphthene	9700	סט כ	9700
2,4-Dinitrophenol	24000	OD C	24000
4-Nitrophenol	24000	סט כ	24000
Dibenzofuran	9700	OD C	9700
2,4-Dinitrotoluene	9700	OD (9700
Diethylphthalate	9700	סט כ	9700
4-Chlorophenyl-phenylether) UD	9700
Fluorene	9700	OD C	9700
4-Nitroaniline	24000		24000
4,6-Dinitro-2-methylphenol	24000		24000
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Client ID: S3
Matrix: SOIL

Lab ID (HSN): 15278 DL Filename: 4067L17

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine 4-Bromophenyl-phenylether	9700 UD 9700 UD	9700 9 7 00
Hexachlorobenzene	9700 UD	9700
Pentachlorophenol	24000 UD	24000
Phenanthrene	1000 JD	9700
Anthracene	9 7 00 UD	9700
Carbazole	9 7 00 UD	9700
Di-n-butylphthalate	1600 JDB	9700
Fluoranthene	2700 JD	9700
Pyrene	4100 JD	9700
Butylbenzylphthalate	9700 UD	9700
3,3'-Dichlorobenzidine	9700 UD	9700
Benz(a)anthracene	2300 JD	9700
Chrysene	2400 JD	9700
bis(2-Ethylhexyl)phthalate	6200 JD	9700
Di-n-octylphthalate	9700 UD	9700
Benzo(b) fluoranthene	4200 JD	9700
Benzo(k) fluoranthene	1300 JD	9700
Benzo(a)pyrene	2800 JD	9700
Indeno(1,2,3-cd)pyrene	3000 JD	9700
Dibenz(a,h)anthracene	9700 UD	9700
Benzo(g,h,i)perylene	2900 JD	9700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	36 % JD	25-121%
Phenol-d5	50 % JD	24-113%
2-Chlorophenol-d4	46%JD	20-130%
1,2-Dichlorobenzene-d4	41%JD	20-130%
Nitrobenzene-d5	46% JD	23-120%
2-Fluorobiphenyl	75 % JD	30-115%
2,4,6-Tribromophenol	5 1% JD	19-122%
Terphenyl-d14	110%JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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Client ID: S4 Lab ID (HSN): 15279
Matrix: SOIL Filename: 4068K07

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 1500 uL
Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 5.88

Compounds:	ug/Kg (PPB)	EQL
Phenol	2100 U	2100
bis(2-Chloroethyl)ether	2100 U	2100
2-Chlorophenol	2100 U	2100
1,3-Dichlorobenzene	2100 U	2100
1,4-Dichlorobenzene	2100 U	2100
1,2-Dichlorobenzene	2100 U	2100
2-Methylphenol	2100 U	2100
2,2'-oxybis(1-Chloropropane)	2100 U	2100
4-Methylphenol	2100 U	2100
N-Nitroso-di-n-propylamine	2100 U.	2100
Hexachloroethane	2100 U	2100
Nitrobenzene	2100 U	2100
Isophorone	2100 U	2100
2-Nitrophenol	2100 U	2100
2,4-Dimethylphenol	2100 U	2100
bis(2-Chloroethoxy)methane	2100 U	2100
2,4-Dichlorophenol	2100 U	2100
1,2,4-Trichlorobenzene	2100 U	2100
Naphthalene	2100 U	2100
4-Chloroaniline	2100 U	2100
Hexachlorobutadiene	2100 U	2100
4-Chloro-3-methylphenol	2100 U	2100
2-Methylnaphthalene	2100 U	2100
Hexachlorocyclopentadiene	2100 U	2100
2,4,6-Trichlorophenol	2100 U	2100
2,4,5-Trichlorophenol	5200 U	5200
2-Chloronaphthalene	2100 U	2100
2-Nitroaniline	5200 U	5200
Dimethylphthalate	2100 U	, 2100
Acenaphthylene	2100 U	2100
2,6-Dinitrotoluene	2100 U	2100
3-Nitroaniline	5200 U	5200
Acenaphthene	2100 U	2100
2,4-Difficiophenoi	5200 0	5200
4-Nitrophenol	5200 U	5200
Dibenzofuran	2100 U	2100
2,4-Dinitrotoluene	2100 U	2100
Diethylphthalate	2100 U	2100
4-Chlorophenyl-phenylether	2100 U	2100
Fluorene	2100 U	2100
4-Nitroaniline	5200 U	5200
4,6-Dinitro-2-methylphenol	5200 U	5200

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Client ID: S4 Lab ID (HSN): 15279
Matrix: SOIL Filename: 4068K07

N-Nitrosodiphenylamine 2100 U 2100 4-Bromophenyl-phenylether 2100 U 2100 Hexachlorobenzene 2100 U 2100 Pentachlorophenol 5200 U 5200 Phenanthrene 1800 J 2100 Anthracene 240 J 2100 Carbazole 2100 U 2100 Di-n-butylphthalate 2100 U 2100 Fluoranthene 2100 U 2100 Pyrene 2100 U 2100 Butylbenzylphthalate 2100 U Y 2100 Benz (a) anthracene 2100 U Y 2100 Chrysene 2100 U Y 2100 Di-n-octylphthalate 6300 T Y 2100 Benzo (b) fluoranthene 4200 U 2100 Benzo (k) fluoranthene 1300 J 2100 Benzo (a) pyrene 2500 U 2100 Indeno (1, 2, 3-cd) pyrene 2600 U 2100	Compounds:	ug/Kg (PPB)	EQL
Dibenz(a,h)anthracene 2100 U 2100	N-Nitrosodiphenylamine 4-Bromophenyl-phenylether Hexachlorobenzene Pentachlorophenol Phenanthrene Anthracene Carbazole Di-n-butylphthalate Fluoranthene Pyrene Butylbenzylphthalate 3,3'-Dichlorobenzidine Benz (a) anthracene Chrysene bis (2-Ethylhexyl) phthalate Di-n-octylphthalate Benzo (b) fluoranthene Benzo (c) pyrene	2100 U 2100 U 2100 U 5200 U 1800 J 240 J 2100 U 2100 U 2100 U 2100 U Y 2100 U Y	2100 2100 2100 5200 2100 2100 2100 2100

Surrogate Recovery		QC LIMITS
2-Fluorophenol	24%J	25-121%
Phenol-d5	48%	24-113%
2-Chlorophenol-d4	46%	20-130%
1,2-Dichlorobenzene-d4	50 % J	20-130%
Nitrobenzene-d5	42 % J	23-120%
2-Fluorobiphenyl	107%	30-115%
2,4,6-Tribromophenol	52%	19-122%
Terphenyl-d14	115%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

3 = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846, November 1986, 3rd Edition.

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Client ID: S4 Lab ID (HSN): 15279 DL Matrix: SOIL Filename: 4067L14

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 1500 uL
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 5.88

Compounds:	ug/Kg (PPB)	EQL
Phenol	21000 UD	21000
bis(2-Chloroethyl)ether	21000 UD	21000
2-Chlorophenol	21000 UD	21000
1,3-Dichlorobenzene	21000 UD	21000
1,4-Dichlorobenzene	,21000 UD	21000
1,2-Dichlorobenzene	21000 UD	21000
2-Methylphenol	21000 UD	21000
2,2'-oxybis(1-Chloropropane)	21000 UD	21000
4-Methylphenol	21000 UD	21000
N-Nitroso-di-n-propylamine	21000 UD	21000
Hexachloroethane	21000 UD	21000
Nitrobenzene .,	21000 UD	21000
Isophorone	21000 UD	21000
2-Nitrophenol	21000 UD	21000
2,4-Dimethylphenol	21000 UD	21000
bis(2-Chloroethoxy)methane	21000 UD	21000
2,4-Dichlorophenol	21000 UD	21000
1,2,4-Trichlorobenzene	21000 UD	21000
Naphthalene	21000 UD	21000
4-Chloroaniline	21000 UD	21000
Hexachlorobutadiene	21000 UD	21000
4-Chloro-3-methylphenol	21000 UD	21000
2-Methylnaphthalene	21000 UD	21000
Hexachlorocyclopentadiene	21000 UD	21000
2,4,6-Trichlorophenol	21000 UD	21000
2,4,5-Trichlorophenol	52000 UD	52000
2-Chloronaphthalene	21000 UD	21000
2-Nitroaniline	52000 UD	52000
Dimethylphthalate	21000 UD	21000
Acenaphthylene	21000 UD	21000
2,6-Dinitrotoluene	21000 UD	21000
3-Nitroaniline	52000 UD	52000
Acenaphth ene	21000 UD	21000
2,4-Dinitrophenol	52000 UD	52000
4-Nitrophenol	52000 UD	52000
Dibenzofuran	21000 UD	21000
2,4-Dinitrotoluene	21000 UD	21000
Diethylphthalate	21000 UD	21000
4-Chlorophenyl-phenylether	21000 UD	21000
Fluorene	21000 UD	21000
4-Nitroaniline	52000 UD	52000
4,6-Dinitro-2-methylphenol	52000 UD	52000

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Client ID: S4 Lab ID (HSN): 15279 DL Matrix: SOIL Filename: 4067L14

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	21000 UD	21000
4-Bromophenyl-phenylether	21000 UD	21000
Hexachlorobenzene	21000 UD	21000
Pentachlorophenol	52000 UD	52000
Phenanthrene	2400 JD	21000
Anthracene	21000 UD	21000
Carbazole	21000 UD	21000
Di-n-butylphthalate	21000 UD	21000
Fluoranthene	3800 JD	21000
Pyrene	7300 JD	21000
Butylbenzylphthalate	21000 UD	21000
3,3'-Dichlorobenzidine	21000 UD	21000
Benz(a)anthracene	2500 JD	21000
Chrysene	2800 JD	21000
bis(2-Ethylhexyl)phthalate	14000 JD	21000
Di-n-octylphthalate	, 21000 UD	21000
Benzo(b)fluoranthene	4000 JD	21000
Benzo(k)fluoranthene	21000 UD	21000
Benzo(a)pyrene	2400 JD	21000
Indeno(1,2,3-cd)pyrene	21000 UD	21000
Dibenz(a,h)anthracene	21000 UD	21000
Benzo(g,h,i)perylene	21000 UD	21000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	28 % JD	25-121%
Phenol-d5	62 % JD	24-113%
2-Chlorophenol-d4	52 % JD	20-130%
1,2-Dichlorobenzene-d4	54%JD	20-130%
Nitrobenzene-d5	5 7%J D	23-120%
2-Fluorobiphenyl	102%JD	30-115%
2,4,6-Tribromophenol	5 4%J D	19-122%
Terphenyl-d14	126%JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S4 MS Lab ID (HSN): 15279MS
Matrix: SOIL Filename: 4068K05

Date Sampled: 02/22/94 Sample Size: 15.6 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL
Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 5.88

Compounds:	ug/Kg (PPB)	EQ
Phenol	180 J	140
bis(2-Chloroethyl)ether	1 4 00 U	1400
2-Chlorophenol	1400 U	1400
1,3-Dichlorobenzene	1400 U	1400
1,4-Dichlorobenzene	240 J	1400
1,2-Dichlorobenzene	1400 U	140
2-Methylphenol	1400 U	1400
2,2'-oxybis(1-Chloropropane)	1400 U	1400
4-Methylphenol	1400 U	1400
N-Nitroso-di-n-propylamine	290 J ·	140
Hexachloroethane	1400 U	1400
Nitrobenzene	1400 U	1400
Isophorone	1400 U	1400
2-Nitrophenol	1400 U	1400
2,4-Dimethylphenol	1400 U	140
bis(2-Chloroethoxy)methane	1400 U	140
2,4-Dichlorophenol	1400 U	140
1,2,4-Trichlorobenzene	290 J	140
Naphthalene	1400 U	140
4-Chloroaniline	1400 U	140
Hexachlorobutadiene	1400 U	140
4-Chloro-3-methylphenol	1400 U	140
2-Methylnaphthalene	1400 U	140
Hexachlorocyclopentadiene	1400 U	140
2,4,6-Trichlorophenol	1400 U	1400
2,4,5-Trichlorophenol	3400 U	3400
2-Chloronaphthalene	1400 U	140
2-Nitroaniline	3400 U	3400
Dimethylphthalate	1400 U	. 140
Acenaphthylene	1400 U	1400
2,6-Dinitrotoluene	1400 U	1400
3-Nitro aniline	3400 U	340
Acenaphthene	340 J	140
2,4-Dinitrophenol	. 3400 U	340
4-Nitrophenol	380 J	340
Dibenzofuran	1400 U	1400
2,4-Dinitrotoluene	320 J	140
Diethylphthalate	1400 U	140
4-Chlorophenyl-phenylether	1400 U	140
Fluorene	1400 U	140
4-Nitroaniline	3400 U	340
4,6-Dinitro-2-methylphenol	3400 U	340

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Client ID: S4 MS Lab ID (HSN): 15279MS
Matrix: SOIL Filename: 4068K05

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	1400 U	1400
4-Bromophenyl-phenylether	1400 U	1400
Hexachlorobenzene	1400 U	1400
Pentachlorophenol	3400 U	3400
Phenanthrene	650 J	1400
Anthracene	150 J	1400
Carbazole	1400 U	1400
Di-n-butylphthalate	1400 U	
Fluoranthene	640 J	1400
Pyrene	1300 J	
Butylbenzylphthalate	1400 U	1400
3,3'-Dichlorobenzidine	1400 U	
Benz(a)anthracene	370 J	1400
Chrysene	400 J	1400
bis(2-Ethylhexyl)phthalate	1200 J	1400
Di-n-octylphthalate	1400 U	1400
Benzo(b) fluoranthene	480 J	1400
Benzo(k) fluoranthene	160 J	1400
Benzo(a)pyrene	310 J	1400
Indeno(1,2,3-cd)pyrene	330 J	1400
Dibenz (a, h) anthracene	1400 U	1400
Benzo(g,h,i)perylene	1400 U	1400

Surrogate Recovery		QC LIMITS
2-Fluorophenol	1 % J	25-121%
Phenol-d5	3 % J	24-113%
2-Chlorophenol-d4	3 % J	20-130%
1,2-Dichlorobenzene-d4	7 % J	20-130%
Nitrobenzene-d5	8 % J	23-120%
2-Fluorobiphenyl	9 % J	30-115%
2,4,6-Tribromophenol	0 % U	19-122%
Terphenyl-414	10 % ፓ	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLMO1)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

Client ID: S4 MSD Lab ID (HSN): 15279MSD Matrix: SOIL Filename: 4068K06

Date Sampled: 02/22/94 Sample Size: 15.3 grams
Date Received: 02/24/94 Extract Vol.: 2000 uL

Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 5.88

Compounds:	ug/Kg (PP	B) EQL
Phenol	6100	2800
bis(2-Chloroethyl)ether	2800	
2-Chlorophenol	2500	
1,3-Dichlorobenzene	2800	
1,4-Dichlorobenzene	1700	
1,2-Dichlorobenzene	2800	
2-Methylphenol	2800	
2,2'-oxybis(1-Chloropropane)	2800	
4-Methylphenol	2800	
N-Nitroso-di-n-propylamine	2100	
Hexachloroethane	2800	
Nitrobenzene	2800	
Isophorone	2800	
2-Nitrophenol	2800	
2,4-Dimethylphenol	2800	
bis(2-Chloroethoxy)methane	2800	
2,4-Dichlorophenol	2800	
1,2,4-Trichlorobenzene	2800	2800
Naphthalene	2800	
4-Chloroaniline	2800	
Hexachlorobutadiene	2800	
4-Chloro-3-methylphenol	2900	2800
2-Methylnaphthalene	2800	U 2800
Hexachlorocyclopentadiene	2800	U 2800
2,4,6-Trichlorophenol	2800	U 2800
2,4,5-Trichlorophenol	6900	U 6900
2-Chloronaphthalene	2800	Ŭ 2800
2-Nitroaniline	6900	U 6900
Dimethylphthalate	2800	U 2800
Acenaphth ylene	2800	U 2800
2,6-Dinitrotoluene	2800	U 2800
3-Nitroan iiine	6900	
Acenaphthene	2900	2800
2,4-Dinitrophenol	6900	
4-Nitrophenol	6900	
Dibenzofuran	2800	
2,4-Dinitrotoluene	3100	2800
Diethylphthalate	2800	
4-Chlorophenyl-phenylether	2800	
Fluorene	2800	
4-Nitroaniline	6900	
4,6-Dinitro-2-methylphenol	6900	ឋ 6900

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Client ID: S4 MSD Lab ID (HSN): 15279MSD Matrix: SOIL Filename: 4068K06

ug/Kg (PPB)	EQL
2800 U	2800
2800 U	2800
2800 U	2800
6900 U	6900
1600 J	2800
2800 U	2800
2800 U	2800
17000 B	2800
2800 U	2800
6900 J Y	2800
2800 U Y	2800
2800 U Y	2800
2400 J Y	2800
3000 7 Y	2800
6500 🐔 Y	2800
2800 U	2800
4800	2800
1600 J	2800
2600 J	2800
2800	2800
2800 U	2800
2900	2800
	2800 U 2800 U 2800 U 6900 U 1600 J 2800 U 17000 B 2800 U 6900 T Y 2800 U Y 2800 U Y 2800 U Y 2400 J Y 3000 T Y 6500 T Y 2800 U 4800 U 1600 J 2800 U

Surrogate Recovery		QC LIMITS
2-Fluorophenol	33 % J	25-121%
Phenol-d5	5 2% J	24-113%
2-Chlorophenol-d4	48 % J	20-130%
1,2-Dichlorobenzene-d4	5 4% J	20-130%
Nitrobenzene-d5	58 % J	23-120%
2-Fluorobiphenyl	94%	30-115%
2,4,6-Tribromophenol	38 % J	19-122%
Terphenyl-d14	7 7 %J	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846, November 1986, 3rd Edition.



Client ID: S5 Lab ID (HSN): 15283 Matrix: SOIL Filename: 4068K11

Date Analyzed: 03/09/94

Date Sampled: 02/22/94
Date Received: 02/24/94
Date Extracted: 03/03/94 Sample Size: 15.4 grams
Extract Vol.: 1000 uL Dil. Factor: GPC Factor: 1

2 % Moisture: 22.42

Compounds:	ug/Kg (PPB)	EQL
Phenol	190 Ј	1700
bis(2-Chloroethyl)ether	1700 U	1700
2-Chlorophenol	1700 U	1700
1,3-Dichlorobenzene	1700 U	1700
1,4-Dichlorobenzene	1700 U	1700
1,2-Dichlorobenzene	1700 U	1700
2-Methylphenol	1700 U	1700
2,2'-oxybis(1-Chloropropane)	1700 U	1700
4-Methylphenol	1700 U	1700
N-Nitroso-di-n-propylamine	1700 U	1700
Hexachloroethane	1700 U	1700
Nitrobenzene	′- 1700 Ŭ	1700
Isophorone	1700 U	1700
2-Nitrophenol	1700 U	1700
2,4-Dimethylphenol	1700 U	1700
bis(2-Chloroethoxy)methane	1700 U	1700
2,4-Dichlorophenol	1700 U	1700
1,2,4-Trichlorobenzene	720 J	1700
Naphthalene	220 J	1700
4-Chloroaniline	1700 U	1700
Hexachlorobutadiene	1700 U	1700
4-Chloro-3-methylphenol	1700 U	1700
2-Methylnaphthalene	1700 U	1700
Hexachlorocyclopentadiene	1700 U	1700
2,4,6-Trichlorophenol	1700 U	1700
2,4,5-Trichlorophenol	4200 U	4200
2-Chloronaphthalene	1700 บ	1700
2-Nitroaniline	4200 U	4200
Dimethylphthalate	1700 U	1700
Acenaphthylene	1700 บ	1700
2,6-Dinit rotol uene	1700 U	1700
3-Nitro aniline	4200 U	4200
Acenaphth ene	3 4 0 J	1700
2,4-Dinitrophenol	4200 U	4200
4-Nitrophenol	4200 U	4200
Dibenzofuran	1700 U	1700
2,4-Dinitrotoluene	1700 U	1700
Diethylphthalate	1700 U	1700
4-Chlorophenyl-phenylether	1700 U	1700
Fluorene	410 J	1700
4-Nitroaniline	4200 U	4200
4,6-Dinitro-2-methylphenol	4200 U	4200

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Client ID: S5 Lab ID (HSN): 15283
Matrix: SOIL Filename: 4068K11

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Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	1700 บ	1700
4-Bromophenyl-phenylether	1700 U	1700
Hexachlorobenzene	1700 U	1700
Pentachlorophenol	4200 U	4200
Phenanthrene	3500	1700
Anthracene	830 J	1700
Carbazole	1700 U	1700
Di-n-butylphthalate	1300 JB	1700
Fluoranthene	3900	1700
Pyrene	7900 T Y	1700
Butylbenzylphthalate	1700 U Y	1700
3,3'-Dichlorobenzidine	1700 U Y	1700
Benz(a)anthracene	3800 🔨 Y	1700
Chrysene	3800 🏅 Y	1700
bis(2-Ethylhexyl)phthalate	30000 E Y	1700
Di-n-octylphthalate	1 7 00 U	1700
Benzo(b) fluoranthene	5800	1700
Benzo(k)fluoranthene	2200	1700
Benzo(a)pyrene	4400	1700
Indeno(1,2,3-cd)pyrene	4200	1700
Dibenz(a,h)anthracene	1700 U	1700
Benzo(g,h,i)perylene	4300	1700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	30 %	25-121%
Phenol-d5	37%	24-113%
2-Chlorophenol-d4	40%	20-130%
1,2-Dichlorobenzene-d4	46%	20-130%
Nitrobenzene-d5	31 % J	23-120%
2-Fluorobinhenyl	72%	30-115%
2,4,6-Triligemophenol	32%	19-122%
Terphenyl-114	77%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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Client ID: S5 Lab ID (HSN): 15283 DL Matrix: SOIL Filename: 4067L18

Date Sampled: 02/22/94 Sample Size: 15.4 grams Date Received: 02/24/94 Extract Vol.: 1000 uL Date Extracted: 03/03/94 Dil. Factor: 10 Date Analyzed: 03/09/94 GPC Factor: 2 % Moisture:

Compounds: ug/Kg (PPB) EQL Phenol 17000 UD 17000 bis(2-Chloroethyl)ether 17000 UD 17000 2-Chlorophenol 17000 UD 17000 1,3-Dichlorobenzene 17000 UD 17000 17000 UD 1,4-Dichlorobenzene 17000 1,2-Dichlorobenzene 17000 UD 17000 2-Methylphenol 17000 UD 17000 17000 UD 17000 2,2'-oxybis(1-Chloropropane) 17000 UD 4-Methylphenol 17000 N-Nitroso-di-n-propylamine 17000 UD 17000 17000 UD Hexachloroethane 17000 Nitrobenzene 17000 UD 17000 17000 UD Isophorone 17000 17000 UD 2-Nitrophenol 17000 17000 UD 2,4-Dimethylphenol 17000 bis(2-Chloroethoxy)methane 17000 UD 17000 17000 UD 17000 2,4-Dichlorophenol 1,2,4-Trichlorobenzene 17000 UD 17000 17000 UD Naphthalene 17000 4-Chloroaniline 17000 UD 17000 17000 UD Hexachlorobutadiene 17000 4-Chloro-3-methylphenol 17000 UD 17000 17000 UD 2-Methylnaphthalene 17000 Hexachlorocyclopentadiene 17000 UD 17000 2,4,6-Trichlorophenol 17000 UD 17000 2,4,5-Trichlorophenol 42000 UD 42000 17000 UD 2-Chloronaphthalene 17000 42000 UD 2-Nitroaniline 42000 Dimethylphthalate 17000 UD .17000 17000 UD Acenaphthylene 17000 2,6-Dinitrotoluene 17000 UD 17000 3-Nitroaniline 42000 UD 42000 Acenaphthene 17000 UD 17000 42000 UD 42000 2,4-Dinitrophenol 42000 UD 4-Nitrophenol 42000 17000 UD Dibenzofuran 17000 17000 UD 17000 2,4-Dinitrotoluene 17000 UD 17000 Diethylphthalate 17000 UD 17000 4-Chlorophenyl-phenylether Fluorene 17000 UD 17000 42000 UD 4-Nitroaniline 42000 4,6-Dinitro-2-methylphenol 42000 UD 42000

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HPN: 2757

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Client ID: S5 Lab ID (HSN): 15283 DL Matrix: SOIL Filename: 4067L18

Compounds:	ug/Kg (PF	PB)	EQL
N-Nitrosodiphenylamine	17000	UD	17000
4-Bromophenyl-phenylether	17000	UD	17000
Hexachlorobenzene	17000	UD	17000
Pentachlorophenol	42000	UD .	42000
Phenanthrene	4500	JD	17000
Anthracene	17000	UD	17000
Carbazole	17000	UD .	17000
Di-n-butylphthalate	2800	JDB	17000
Fluoranthene	5800	JD	17000
Pyrene	10000	JD	17000
Butylbenzylphthalate	17000	UD	17000
3,3'-Dichlorobenzidine	17000	UD	17000
Benz(a)anthracene	4100		17000
Chrysene	3800	JD	17000
bis(2-Ethylhexyl)phthalate	56000		17000
Di-n-octylphthalate	17000	UD .	17000
Benzo(b) fluoranthene		JD	1 70 00
Benzo(k) fluoranthene		UD	17000
Benzo(a)pyrene		JD	17000
Indeno(1,2,3-cd)pyrene		JD	17000
Dibenz(a,h)anthracene		UD	17000
Benzo(g,h,i)perylene	5000	JD	17000

	QC LIMITS
41%J D	25-121%
49 % JD	24-113%
47% JD	20-130%
46 % JD	20-130%
48% JD	23-120%
6 6% JD	30-115%
36 % JD	19-122%
96 % JD	18-137%
	49%JD 47%JD 46%JD 48%JD 66%JD 36%JD

TCL = Target Coumpound List EPA Contract Laboratory Program (OLMO1)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S6 Lab ID (HSN): 15287 Matrix: SOIL Filename: 4068K12

Date Sampled: 02/22/94 Sample Size: 15.2 grams Date Received: 02/24/94 Extract Vol.: 1000 uL Date Extracted: 03/03/94 Dil. Factor: 1 Date Extracted: 03/03/94
Date Analyzed: 03/10/94

GPC Factor: 2 % Moisture: 17.96

ompounds:	ug/Kg (PPB)	EQ
Phenol	1600 U	160
bis(2-Chloroethyl)ether	1600 U	160
2-Chlorophenol	1600 U	160
1,3-Dichlorobenzene	1600 U	160
1,4-Dichlorobenzene	1600 U	160
1,2-Dichlorobenzene	1600 U	160
2-Methylphenol	1600 U	160
2,2'-oxybis(1-Chloropropane)	1600 U	160
4-Methylphenol	1600 U	160
N-Nitroso-di-n-propylamine	1600 U	160
Hexachloroethane	1600 U	160
Nitrobenzene	′~ 1600 U	160
Isophorone	1600 U	160
2-Nitrophenol	1600 U	160
2,4-Dimethylphenol	1600 U	160
bis(2-Chloroethoxy)methane	1600 U	160
2,4-Dichlorophenol	1600 U	160
1,2,4-Trichlorobenzene	690 J	160
Naphthalene	510 J	160
4-Chloroaniline	1600 U	160
Hexachlorobutadiene	1600 U	160
4-Chloro-3-methylphenol	1600 U	160
2-Methylnaphthalene	330 J	160
Hexachlorocyclopentadiene	1600 U	160
2,4,6-Trichlorophenol	1600 U	160
2,4,5-Trichlorophenol	4000 U	400
2-Chloronaphthalene	1600 U	160
2-Nitroaniline	4000 U	400
Dimethylphthalate	1600 U	160
Acenaphth ylene	1600 U	160
2,6-Dinitrotoluene	1600 U	160
3-Nitroaniline	4000 U	400
Acenaphth ene	740 J	160
2,4-Dinitrophenol	4000 U	400
4-Nitrophenol	4000 U	400
Dibenzofuran	760 J	160
2,4-Dinitrotoluene	1600 U	160
Diethylphthalate	1600 U	160
4-Chlorophenyl-phenylether	1600 U	160
Fluorene	1100 J	160
4-Nitroaniline	4000 U	400
4,6-Dinitro-2-methylphenol	4000 U	400

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Client ID: S6 Lab ID (HSN): 15287
Matrix: SOIL Filename: 4068K12

EQL
500
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500

Surrogate Recovery		QC LIMITS
2-Fluorophenol	45%	25-121%
Phenol-d5	47%	24-113%
2-Chlorophenol-d4	54%	20-130%
1,2-Dichlorobenzene-d4	67%	20-130%
Nitrobenzene-d5	43%	23-120%
2-Fluorobiphenyl	77%	30-115%
2,4,6-Tribromophenol	31%	19-122%
Terphenyl-d14	82%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S6 Lab ID (HSN): 15287 DL Matrix: SOIL Filename: 4067L19

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 17.96

Compounds:	ug/Kg (PPB)	EQL
Phenol	16000 UD	16000
bis(2-Chloroethyl)ether	16000 UD	16000
2-Chlorophenol	16000 UD	16000
1,3-Dichlorobenzene	16000 UD	16000
1,4-Dichlorobenzene	16000 UD	16000
1,2-Dichlorobenzene	16000 UD	16000
2-Methylphenol	16000 UD	16000
2,2'-oxybis(1-Chloropropane)	16000 UD	16000
4-Methylphenol	16000 UD	16000
N-Nitroso-di-n-propylamine	16000 UD	16000
Hexachloroethane	16000 UD	16000
Nitrobenzene	16000 UD	16000
Isophorone	16000 UD	16000
2-Nitrophenol	16000 UD	16000
2,4-Dimethylphenol	16000 UD	16000
bis(2-Chloroethoxy)methane	16000 UD	16000
2,4-Dichlorophenol	16000 UD	16000
1,2,4-Trichlorobenzene	16000 UD	16000
Naphthalene	16000 UD	16000
4-Chloroaniline	16000 UD	16000
Hexachlorobutadiene	16000 UD	16000
4-Chloro-3-methylphenol	16000 UD	16000
2-Methylnaphthalene	16000 UD	16000
Hexachlorocyclopentadiene	16000 UD	16000
2,4,6-Trichlorophenol	16000 VD	16000
2,4,5-Trichlorophenol	40000 UD	40000
2-Chloronaphthalene	16000 UD	16000
2-Nitroaniline	40000 UD	40000
Dimethylphthalate	16000 UD	16000
Acenaphthylene	16000 UD	16000
2,6-Dinitrotoluene	16000 UD	16000
3-Nitroaniline	40000 UD 16000 UD	40000
Acenaphthene	40000 UD	16000 40000
2,4-Dinitrophenol	40000 UD	40000
4-Nitrophenol Dibenzofuran	16000 UD	16000
2,4-Dinitrotoluene	16000 UD	16000
Diethylphthalate	16000 UD	16000
4-Chlorophenyl-phenylether	16000 UD	16000
Fluorene	16000 UD	16000
4-Nitroaniline	40000 UD	40000
4,6-Dinitro-2-methylphenol	40000 UD	40000
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Client ID: S6 Lab ID (HSN): 15287 DL Matrix: SOIL Filename: 4067L19

Compounds:	ug/Kg (PPB)	EQL
N Nitrogodinhonvlamino	16000 170	
N-Nitrosodiphenylamine	16000 UD	16000
4-Bromophenyl-phenylether	16000 UD	16000
Hexachlorobenzene	16000 UD	16000
Pentachlorophenol	40000 UD	40000
Phenanthrene	9700 JD	16000
Anthracene	2900 JD	16000
Carbazole	16000 UD	16000
Di-n-butylphthalate	1800 JDB	16000
Fluoranthene .,	10000 JD	16000
Pyrene	1 7 000 D	16000
Butylbenzylphthalate	16000 UD	16000
3,3'-Dichlorobenzidine	16000 UD	16000
Benz(a)anthracene	6400 JD	16000
Chrysene	5900 JD	16000
bis(2-Ethylhexyl)phthalate	42000 D	16000
Di-n-octylphthalate	16000 UD	16000
Benzo(b) fluoranthene	8000 JD	16000
Benzo(k) fluoranthene	3500 JD	16000
Benzo(a) pyrene	6000 JD	16000
Indeno(1,2,3-cd)pyrene	5900 JD	16000
Dibenz(a,h)anthracene	16000 UD	16000
Benzo(g,h,i)perylene	5400 JD	16000
penzo (a'u' i) ber liene	2400 00	10000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	63 % JD	25-121%
Phenol-d5	71%JD	24-113%
2-Chlorophenol-d4	65 % JD	20-130%
1,2-Dichlorobenzene-d4	6 6%J D	20-130%
Nitrobenzene-d5	6 6%J D	23-120%
2-Fluorobiphenyl	76 % JD	30-115%
2,4,6-Tribromophenol	35 % JD	19-122%
Terphenyl-d14	108%JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846, November 1986, 3rd Edition.

Huntingdon

Client ID: S7 Lab ID (HSN): 15289
Matrix: SOIL Filename: 4068K13

Date Sampled: 02/22/94 Sample Size: 15.3 grams
Date Received: 02/24/94 Extract Vol.: 500 uL
Date Extracted: 03/03/94 Dil. Factor: 1

Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/10/94 GPC Factor: 2
% Moisture: 30.67

Compounds:	ug/Kg (PPB)	EQL
Phenol	9 4 0 U	940
bis(2-Chloroethyl)ether	9 4 0 U	940
2-Chlorophenol	9 4 0 U	940
1,3-Dichlorobenzene	940 U	940
1,4-Dichlorobenzene	940 U	940
1,2-Dichlorobenzene	9 4 0 U	940
2-Methylphenol	940 U	940
2,2'-oxybis(1-Chloropropane)	940 U	940
4-Methylphenol	940 U	940
N-Nitroso-di-n-propylamine	940 U	940
Hexachloroethane	940 U	940
Nitrobenzene	940 U	940
Isophorone	940 U	940
2-Nitrophenol	940 U	940
2,4-Dimethylphenol	940 U	940
bis (2-Chloroethoxy) methane	940 U	940
2,4-Dichlorophenol	940 U	940
1,2,4-Trichlorobenzene	430 J	940
Naphthalene	260 J	940
4-Chloroaniline	940 U	940
Hexachlorobutadiene	940 U	940
4-Chloro-3-methylphenol	940 U	940
2-Methylnaphthalene Hevachlorocyclopentadiene	550 J 940 U	940 940
Hexachlorocyclopentadiene 2,4,6-Trichlorophenol	940 U 940 U	940
2,4,5-Trichlorophenol	2400 U	2400
2.Chloronaphthalene	2400 U 940 U	2400 940
2-Chioronaphchaiene 2-Nitroaniline	2400 U	2400
Dimethylphthalate	940 U	940
Acenaphthylene	940 U	940
2,6-Dinitrotoluene	940 U	940
3-Nitroaniline	2400 U	2400
Acenaphthene	250 J	940
2,4-Dinitrophenol	· * 2400 U	2400
4-Nitrophenol	2400 U	2400
Dibenzofuran	280 J	940
2,4-Dinitrotoluene	940 U	940
Diethylphthalate	940 U	940
4-Chlorophenyl-phenylether	9 4 0 U	940
Fluorene	290 J	940
4-Nitroaniline	2400 U	2400
4,6-Dinitro-2-methylphenol	2400 U	2400
(continued)		

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Client ID: S7
Matrix: SOIL

Lab ID (HSN): 15289 Filename: 4068K13

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	940 U	940
4-Bromophenyl-phenylether	940 U	940
Hexachlorobenzene	9 4 0 U	940
Pentachlorophenol	2400 U	2400
Phenanthrene	2100	940
Anthracene	580 J	940
Carbazole	9 4 0 U	940
Di-n-butylphthalate	800 JB	940
Fluoranthene	2700	940
Pyrene	5400 T Y	940
Butylbenzylphthalate	940 U Y	940
3,3'-Dichlorobenzidine	940 U Y	940
Benz(a)anthracene	2100 - Y	
Chrysene	2300 ° Y	940
bis(2-Ethylhexyl)phthalate	9400 E Y	940
Di-n-octylphthalate	940 U	940
Benzo (b) Iluoranthene	3600	940
Benzo(k)fluoranthene	1200	940
Benzo(a)pyrene	2200	940
Indeno(1,2,3-cd)pyrene	2100	940
Dibenz(a,h)anthracene	460 J	940
Benzo(g,h,i)perylene	1900	940

Surrogate Recovery		QC LIMITS
2-Fluorophenol	30%	25-121%
Phenol-d5	36%	24-113%
2-Chlorophenol-d4	41%	20-130%
1,2-Dichlorobenzene-d4	41%	20-130%
Nitrobenzene-d5	30%	23-120%
2-Fluorobiphenyl	71%	30-115%
2,4,6-Tribromophenol	47%	19-122%
Terphenyl-d14	75%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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.: "Suiting Engineers of Environmental Scientists"

Client ID: S7 Lab ID (HSN): 15289 DL Matrix: SOIL Filename: 4067L20

Date Sampled: 02/22/94 Sample Size: 15.3 grams
Date Received: 02/24/94 Extract Vol.: 500 uL
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 30.67

Compounds:	ug/Kg (PPB)	EQL
Phenol	9400 UD	9400
bis(2-Chloroethyl)ether	9 400 UD	9400
2-Chlorophenol	9400 UD	9400
1,3-Dichlorobenzene	9400 UD	9400
1,4-Dichlorobenzene	9400 UD	9400
1,2-Dichlorobenzene	9400 UD	9400
2-Methylphenol	9400 UD	9400
2,2'-oxybis(1-Chloropropane)	9400 UD	9400
4-Methylphenol	9400 UD	9400
N-Nitroso-di-n-propylamine	9400 UD	9400
Hexachloroethane	9400 UD	9400
Nitrobenzene	9400 UD	9400
Isophorone	9400 UD	9400
2-Nitrophenol	9400 UD	9400
2,4-Dimethylphenol	9400 UD	9400
bis(2-Chloroethoxy)methane	9400 UD	9400
2,4-Dichlorophenol	9400 UD	9400
1,2,4-Trichlorobenzene	9400 UD	9400
Naphthalene	9400 UD	9400
4-Chloroaniline	9400 UD	9400
Hexachlorobutadiene	9400 UD	9400
4-Chloro-3-methylphenol	9400 UD	9400
2-Methylnaphthalene	9400 UD	9400
Hexachlorocyclopentadiene	9400 UD	9400
2,4,6-Trichlorophenol	9400 UD	9400
2,4,5-Trichlorophenol	24000 UD	24000
2-Chloronaphthalene	9400 UD	9400
2-Nitroaniline	24000 UD	24000
Dimethylphthalate	9400 UD	9400
Acenaphthylene	9400 UD	9400
2,6-Dinitrotoluene	9400 UD	9400
3-Nitroaniline	24000 UD	24000
Acenaphthene	9400 UD	9400
2,4-Dinitrophenol	24000 UD	24000
4-Nitrophenol	24000 UD	24000
Dibenzofuran	9400 UD	9400
2,4-Dinitrotoluene	9400 UD	9400
Diethylphthalate	9400 UD	9400
4-Chlorophenyl-phenylether	9400 UD	9400
Fluorene	9400 UD	9400
4-Nitroaniline	24000 UD	24000
4,6-Dinitro-2-methylphenol	24000 UD	24000

(continued)

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Client ID: S7 Lab ID (HSN): 15289 DL Matrix: SOIL Filename: 4067L20

			
Compounds:	ug/Kg (Pl	PB)	EQL
N-Nitrosodiphenylamine	9400	UD	9400
4-Bromophenyl-phenylether	9400	WD	9400
Hexachlorobenzene	9400	UD	9400
Pentachlorophenol	24000	UD	24000
Phenanthrene	2200	JD	9400
Anthracene	9400	UD	9400
Carbazole	9400	UD	9400
Di-n-butylphthalate	1100	JDB	9400
Fluoranthene	3200	JD	9400
Pyrene	7700	JD	9400
Butylbenzylphthalate	9400	UD	9400
3,3'-Dichlorobenzidine	9400	UD	9400
Benz(a)anthracene	2200	JD	9400
Chrysene	2400	JD	9400
bis(2-Ethylhexyl)phthalate	18000	D	9400
Di-n-octylphthalate	9400	UD	9400
Benzo(b) fluoranthene	3700	JD	9400
Benzo(k) fluoranthene	940	JD	9400
Benzo(a)pyrene	2400	JD	9400
Indeno(1,2,3-cd)pyrene	2400	JD	9400
Dibenz(a,h)anthracene	9400	UD	9400
Benzo(g,h,i)perylene	2500	JD	9400

Surrogate Recovery		QC LIMITS
2-Fluorophenol	39 % JD	25-121%
Phenol-d5	51 % JD	24-113%
2-Chlorophenol-d4	46 % JD	20-130%
1,2-Dichlorobenzene-d4	39 % JD	20-130%
Nitrobenzene-d5	45 % JD	23-120%
2-Fluorobiphenyl	70 % JD	30-11 5%
2,4,6-Tribromophenol	55 % JD	19-122%
Terphenyl-dl4	108 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S8 Lab ID (HSN): 15290 Matrix: SOIL Filename: 4068K14

Sample Size: 15.3 grams 1000 uL Date Sampled: 02/22/94 Extract Vol.: Dil. Factor: Date Received: 02/24/94 Date Extracted: 03/03/94 Date Extracted: 03/03/94
Date Analyzed: 03/10/94 1

GPC Factor: 2 % Moisture: 23.54

Compounds:	ug/Kg (PPB)	EQ
Phenol	1700 U	170
bis(2-Chloroethyl)ether	1700 U	170
2-Chlorophenol	1700 ป	170
1,3-Dichlorobenzene	1700 U	170
1,4-Dichlorobenzene	1700 U	170
1,2-Dichlorobenzene	1700 U	170
2-Methylphenol	1700 U	170
2,2'-oxybis(1-Chloropropane)	1700 U	170
4-Methylphenol	1700 ับ	170
N-Nitroso-di-n-propylamine	1700 U .	170
Hexachloroethane	1700 U	170
Nitrobenzene	1700 U	170
Isophorone	1 7 00 U	170
2-Nitrophenol	1700 U	170
2,4-Dimethylphenol	1700 U	170
bis(2-Chloroethoxy)methane	1700 U	170
2,4-Dichlorophenol	1700 U	170
1,2,4-Trichlorobenzene	1700 U	170
Naphthalene	250 J	170
4-Chloroaniline	1700 U	170
Hexachlorobutadiene	1700 U	170
4-Chloro-3-methylphenol	1700 U	170
2-Methylnaphthalene	480 J	170
Hexachlorocyclopentadiene	1700 U	170
2,4,6-Trichlorophenol	1700 U	170
2,4,5-Trichlorophenol	4300 U	430
2-Chloronaphthalene	1700 U	170
2-Nitroaniline	4300 U	430
Dimethylphthalate	1700 U	· 170
Acenaphthylene	1700 U	170
2,6-Dinitrotoluene	1700 U	170
3-Nitroaniline	4300 U	430
Acenaphth ene	520 J	170
2,4-Dinitrophenol	4300 U	430
4-Nitrophenol	4300 U	430
Dibenzofuran	230 J	170
2,4-Dinitrotoluene	1700 U	170
Diethylphthalate	1700 U	170
4-Chlorophenyl-phenylether	1700 U	170
Fluorene	200 Ј	170
4-Nitroaniline	4300 U	430
4,6-Dinitro-2-methylphenol	4300 U	430

(continued)

Client ID: S8
Matrix: SOIL

Lab ID (HSN): 15290 Filename: 4068K14

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	1700 U	1700
4-Bromophenyl-phenylether	1700 U	1700
Hexachlorobenzene	1700 U	1700
Pentachlorophenol	4300 U	4300
Phenanthrene	2900	1700
Anthracene	420 J	1700
Carbazole	1700 U	1700
Di-n-butylphthalate	3900 B	1700
Fluoranthene	3600	1700
Pyrene	7600 ブ Y	1700
Butylbenzylphthalate	4500 T Y	1700
3,3'-Dichlorobenzidine	1700 U Y	
Benz(a)anthracene	3800 T Y	1700
Chrysene	4700 🕻 Y	1700
bis(2-Ethylhexyl)phthalate	4700 ≤ Y	1700
Di-n-octylphthalate	1700 U	1700
Benzo(b)fluoranthene	7600	1700
Benzo(k)fluoranthene	2400	1700
Benzo(a)pyrene	5200	1700
Indeno(1,2,3-cd)pyrene	6300	1700
Dibenz(a,h)anthracene	1100 J	1700
Benzo(g,h,i)perylene	6100	1700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	43%	25-121%
Phenol-d5	45%	24-113%
2-Chlorophenol-d4	55%	20-130%
1,2-Dichlorobenzene-d4	56 %	20-130%
Nitrobenzene-d5	36 % J	23-120%
2-Fluorobiphenyl	75%	30-115%
2,4,6-Tribromophenol	61%	19-122%
Terphenyl-d14	96 %	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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Client ID: S8 Lab ID (HSN): 15290 DL Matrix: SOIL Filename: 4067L21

Date Sampled: 02/22/94 Sample Size: 15.3 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 23.54

Compounds:	ug/Kg (PPB)	EQL
Phenol	17000 UD	17000
bis(2-Chloroethyl)ether	17000 UD	17000
2-Chlorophenol	17000 UD	17000
1,3-Dichlorobenzene	17000 UD	17000
1,4-Dichlorobenzene	17000 UD	17000
1,2-Dichlorobenzene	17000 UD	17000
2-Methylphenol	17000 UD	17000
2,2'-oxybis(1-Chloropropane)	17000 UD	17000
4-Methylphenol	17000 UD	17000
N-Nitroso-di-n-propylamine	17000 UD	17000
Hexachloroethane	17000 UD	17000
Nitrobenzene	17000 UD	17000
Isophorone	17000 UD	17000
2-Nitrophenol	17000 UD	17000
2,4-Dimethylphenol	17000 UD	17000
bis (2-Chloroethoxy) methane	17000 UD	17000
2,4-Dichlorophenol	17000 UD	17000
1,2,4-Trichlorobenzene	17000 UD	17000
Naphthalene	17000 UD	17000
4-Chloroaniline	17000 UD	17000
Hexachlorobutadiene	17000 UD	17000
4-Chloro-3-methylphenol	17000 UD	17000
2-Methylnaphthalene	17000 UD	17000
Hexachlorocyclopentadiene	17000 UD	17000
2,4,6-Trichlorophenol	17000 UD	17000
2,4,5-Trichlorophenol	43000 UD	43000
2-Chloronaphthalene	17000 UD	17000
2-Nitroaniline	43000 UD	43000
Dimethylphthalate	17000 UD	17000
Acenaphthylene	17000 UD	17000
2,6-Dinitrotoluene	17000 UD	17000
3-Nitroaniline	43000 UD	43000
Acenaphthene	17000 UD	17000
2,4-Dinitrophenol	43000 UD	43000
4-Nitrophenol	43000 UD	43000
Dibenzofuran	17000 UD	17000
2,4-Dinitrotoluene	17000 UD	17000
Diethylphthalate	17000 UD	17000
4-Chlorophenyl-phenylether	17000 UD	17000
Fluorene	17000 UD	17000
4-Nitroaniline	43000 UD	43000
4,6-Dinitro-2-methylphenol	43000 UD	43000
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Client ID: S8
Matrix: SOIL

Lab ID (HSN): 15290 DL Filename: 4067L21

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	17000 UD	17000
4-Bromophenyl-phenylether	17000 UD	17000
Hexachlorobenzene	17000 UD	17000
Pentachlorophenol	43000 UD	43000
Phenanthrene	3200 JD	17000
Anthracene	17000 UD	17000
Carbazole	17000 UD	17000
Di-n-butylphthalate	7000 JDB	17000
Fluoranthene	5400 JD	17000
Pyrene	8700 JD	17000
Butylbenzylphthalate	9100 JD	17000
3,3'-Dichlorobenzidine	17000 UD	17000
Benz(a)anthracene	4000 JD	17000
Chrysene	4600 JD	17000
bis(2-Ethylhexyl)phthalate	9300 JD	17000
Di-n-octylphthalate	17000 UD	17000
Benzo(b)fluoranthene	7600 JD	17000
Benzo(k)fluoranthene	2500 JD	17000
Benzo(a)pyrene	4900 JD	17000
Indeno(1,2,3-cd)pyrene	7500 JD	17000
Dibenz(a,h)anthracene	17000 UD	17000
Benzo(g,h,i)perylene	7200 JD	17000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	57 % JD	25-121%
Phenol-d5	65 % JD	24-113%
2-Chlorophenol-d4	61%JD	20-130%
1,2-Dichlorobenzene-d4	5 4%J D	20-130%
Nitrobenzene-d5	56 % JD	23-120%
2-Fluorobiphenyl	. 76 % JD	30-115%
2,4,6-Tribromophenol	61 % JD	19-122%
Terphenyl-d14	108%JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S9 Lab ID (HSN): 15291 Matrix: SOIL Filename: 4068K15

Sample Size: 15.1 grams Extract Vol.: 500 uL

Date Sampled: 02/22/94
Date Received: 02/24/94
Date Extracted: 03/03/94 Dil. Factor: 1 Date Analyzed: 03/10/94 GPC Factor: 2

% Moisture: 14.95

Compounds:	ug/Kg (PPB)	EQL
Phenol	78	10 U	780
bis(2-Chloroethyl)ether		0 U	780
2-Chlorophenol		O U	780
1,3-Dichlorobenzene		0 U	780
1,4-Dichlorobenzene		0 U	780
1,2-Dichlorobenzene		0 U	780
2-Methylphenol		0 U	780
2,2'-oxybis(1-Chloropropane)		U 0	780
4-Methylphenol		υ 0	780
N-Nitroso-di-n-propylamine		U 0	780
Hexachloroethane		0 U	780
Nitrobenzene		10 U	780
Isophorone		0 U	780
2-Nitrophenol		10 U	780
2,4-Dimethylphenol		10 U	780
bis(2-Chloroethoxy)methane		0 U	780
2,4-Dichlorophenol		10 U	780
1,2,4-Trichlorobenzene		.0 J	780
Naphthalene		0 U	780
4-Chloroaniline		0 U	780
Hexachlorobutadiene		0 U	780
4-Chloro-3-methylphenol		0 U	780
2-Methylnaphthalene		0 J	780
Hexachlorocyclopentadiene		υ ο υ	780
2,4,6-Trichlorophenol		0 U	780
2,4,5-Trichlorophenol		0 U	1900
2-Chloronaphthalene		U O	780
2-Nitroaniline		0 U	1900
Dimethylphthalate		O U	780
Acenaphthylene		0 U	780
2,6-Dinitrotoluene		0 U	780
3-Nitroaniline		0 U	1900
Acenaphthene		0 U	780
		0 U	1900
4-Nitrophenol		0 U	1900
Dibenzofuran		10 U	780
2,4-Dinitrotoluene		0 U	780
Diethylphthalate		10 U	780
4-Chlorophenyl-phenylether		10 U	780
Fluorene		.0 J	780
4-Nitroaniline		0 U	1900
4,6-Dinitro-2-methylphenol		υ ο υ	1900
-, o managed a meenjaphener	150		-200
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Client ID: S9 Lab II
Matrix: SOIL Fi

Lab ID (HSN): 15291 Filename: 4068K15

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	780 U	780
4-Bromophenyl-phenylether	780 U	780
Hexachlorobenzene	780 U	780
Pentachlorophenol	1900 ປ	1900
Phenanthrene	1100	780
Anthracene	190 J	780
Carbazole	780 บ	780
Di-n-butylphthalate	620 JB	780
Fluoranthene	1500	780
Pyrene	4100 J Y	780
Butylbenzylphthalate	780 U Y	780
3,3'-Dichlorobenzidine	780 U Y	780
Benz(a)anthracene	1500 T Y	780
Chrysene	1500 5 Y	780
bis(2-Ethylhexyl)phthalate	4300 5 Y	780
Di-n-octylphthalate	780 U	780
Benzo(b) fluoranthene	2800	780
Benzo(k) fluoranthene	860	780
Benzo(a)pyrene	1400	780
Indeno(1,2,3-cd)pyrene	1500	780
Dibenz(a,h)anthracene	780 U	780
Benzo(g,h,i)perylene	1400	780

Surrogate Recovery		QC LIMITS
2-Fluorophenol	30%	25-121%
Phenol-d5	41%	24-113%
2-Chlorophenol-d4	45%	20-130%
1,2-Dichlorobenzene-d4	61%	20-130%
Nitrobenzene-d5	42%	23-120%
2-Fluorobiphenyl	75%	30-115%
2,4,6-Tribromophenol	22%	19-122%
Terphenyl-d14	92 %	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



Client ID: S9 Lab ID (HSN): 15291 DL Matrix: SOIL Filename: 4067L22

Date Sampled: 02/22/94 Sample Size: 15.1 grams
Date Received: 02/24/94 Extract Vol.: 500 uL
Date Extracted: 03/03/94 Dil. Factor: 10

Date Analyzed: 03/09/94 GPC Factor: 2 % Moisture: 14.95

Compounds:	ug/Kg (PPB)	EQL
Phenol	7800 UD	7800
bis(2-Chloroethyl)ether	7800 UD	7800
2-Chlorophenol	7800 UD	7800
1,3-Dichlorobenzene	7800 UD	7800
1,4-Dichlorobenzene	7800 UD	7800
1,2-Dichlorobenzene	7800 UD	7800
2-Methylphenol	7800 UD	7800
2,2'-oxybis(1-Chloropropane)	7800 UD	7800
4-Methylphenol	7800 UD	7800
N-Nitroso-di-n-propylamine	7800 UD	7800
Hexachloroethane	7800 UD	7800
Nitrobenzene	7800 UD	7800
Isophorone	7800 UD	7800
2-Nitrophenol	7800 UD	7800
2,4-Dimethylphenol	7800 UD	7800
bis(2-Chloroethoxy)methane	7800 UD	7800
2,4-Dichlorophenol	7800 UD	7800
1,2,4-Trichlorobenzene	7800 UD	7800
Naphthalene	7800 UD	7800
4-Chloroaniline	7800 UD	7800
Hexachlorobutadiene	7800 UD	7800
4-Chloro-3-methylphenol	7800 UD	7800
2-Methylnaphthalene	7800 UD	7800
Hexachlorocyclopentadiene	7800 UD	7800
2,4,6-Trichlorophenol	7800 UD	7800
2,4,5-Trichlorophenol	19000 UD	19000
2-Chloronaphthalene	7800 UD	7800
2-Nitroaniline	19000 UD	19000
Dimethylphthalate	7800 UD	7800
Acenaphthylene	7800 UD	7800
2,6-Dinitrotoluene	7800 UD	7800
3-Nitroaniline	19000 UD	19000
Acenaphthene	, 7800 UD	7800
2,4-Dinitrophenol	19000 UD	19000
4-Nitrophenol	19000 UD	19000
Dibenzofuran	7800 UD	7800
2,4-Dinitrotoluene	7800 UD	7800
Diethylphthalate	7800 UD	7800
4-Chlorophenyl-phenylether	7800 UD	7800
Fluorene	7800 UD	7800
4-Nitroaniline	19000 UD	19000
4,6-Dinitro-2-methylphenol	19000 UD	19000

(continued)

Huntingdon

Usulina Engineers General Scientists

TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS (CONTINUED) EPA METHOD 8270

Client ID: S9 Lab ID (HSN): 15291 DL Matrix: SOIL Filename: 4067L22

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	7800 UD	7800
4-Bromophenyl-phenylether	7800 UD	7800
Hexachlorobenzene	7800 UD	7800
Pentachlorophenol	19000 UD	19000
Phenanthrene	1400 JD	7800
Anthracene	7800 UD	7800
Carbazole	7800 UD	7800
Di-n-butylphthalate	1300 JDB	7800
Fluoranthene	2600 JD	7800
Pyrene	5200 JD	7800
Butylbenzylphthalate	7800 UD	7800
3,3'-Dichlorobenzidine	7800 UD	7800
Benz(a)anthracene	1600 JD	7800
Chrysene	1900 JD	7800
bis(2-Ethylhexyl)phthalate	9800 D	7800
Di-n-octylphthalate	, 7800 UD	7800
Benzo(b)fluoranthene	3100 JD	7800
Benzo(k) fluoranthene	860 JD	7800
Benzo(a)pyrene	1400 JD	7800
Indeno(1,2,3-cd)pyrene	1800 JD	7800
Dibenz(a,h)anthracene	7800 UD	7800
Benzo(g,h,i)perylene	1700 JD	7800

Surrogate Recovery		QC LIMITS
2-Fluorophenol	39 % JD	25-121%
Phenol-d5	55 % JD	24-113%
2-Chlorophenol-d4	5 3%J D	20-130%
1,2-Dichlorobenzene-d4	58 % JD	20-130%
Nitrobenzene-d5	61 % JD	23-120%
2-Fluorobiphenyl	78 % JD	30-115 %
2,4,6-Tribromophenol	2 4% JD	19-122%
Terphenyl-d14	115 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

= Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.



HPN: 2757

TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Client ID: S10 Lab ID (HSN): 15292 DL Matrix: SOIL Filename: 4067L23

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 2000 uL
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 40.38

Compounds:	ug/Kg (PF	PB)	EQL
Phenol	44000	UD	44000
bis(2-Chloroethyl)ether	44000	UD	44000
2-Chlorophenol	44000	UD	44000
1,3-Dichlorobenzene	44000	UD	44000
1,4-Dichlorobenzene	44000	UD	44000
1,2-Dichlorobenzene	44000	UD	44000
2-Methylphenol	44000	UD	44000
2,2'-oxybis(1-Chloropropane)	44000	UD	44000
4-Methylphenol	44000	UD	44000
N-Nitroso-di-n-propylamine	44000	UD	44000
Hexachloroethane	44000	UD	44000
Nitrobenzene	44000	UD	44000
Isophorone	44000	Œ	44000
2-Nitrophenol	44000	UD	44000
2,4-Dimethylphenol	44000	UD	44000
bis(2-Chloroethoxy)methane	44000	UD	44000
2,4-Dichlorophenol	44000	UD	44000
1,2,4-Trichlorobenzene	44000	UD	44000
Naphthalene	54000	D	44000
4-Chloroaniline	44000	UD	44000
Hexachlorobutadiene	44000		44000
4-Chloro-3-methylphenol	44000	UD	44000
2-Methylnaphthalene	230000	D	44000
Hexachlorocyclopentadiene	44000	UD	44000
2,4,6-Trichlorophenol	44000	UD	44000
2,4,5-Trichlorophenol	110000	Œ	110000
2-Chloronaphthalene	44000	ÚD	44000
2-Nitroaniline	110000	UD	110000
Dimethylphthalate	44000	Œ	. 44000
Acenaphth ylene	44000		44000
2,6-Dinitrotoluene	44000		44000
3-Nitroan iline	110000	Œ	110000
Acenaphth ene	14000		44000
2,4-Dinitrophenol	110000		110000
4-Nitrophenol	110000		110000
Dibenzofuran	44000		44000
2,4-Dinitrotoluene	44000		44000
Diethylphthalate	44000		44000
4-Chlorophenyl-phenylether	44000		44000
Fluorene	32000		44000
4-Nitroaniline	110000	UD	110000
4,6-Dinitro-2-methylphenol	110000	UD	110000

(continued)

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HPN: 2757

TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS (CONTINUED) EPA METHOD 8270

Client ID: S10
Matrix: SOIL

Lab ID (HSN): 15292 DL Filename: 4067L23

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	44000 UD	44000
4-Bromophenyl-phenylether	44000 UD	44000
Hexachlorobenzene	44000 UD	44000
Pentachlorophenol	110000 UD	110000
Phenanthrene	88000 D	44000
Anthracene	7500 JD	44000
Carbazole	44000 UD	44000
Di-n-butylphthalate	44000 UD	44000
Fluoranthene	35000 JD	44000
Pyrene	62000 D	44000
Butylbenzylphthalate	44000 UD	44000
3,3 ⁷ -Dichlorobenzidine	44000 UD	44000
Benz(a)anthracene	20000 JD	44000
Chrysene	20000 JD	44000
bis(2-Ethylhexyl)phthalate	44000 UD	44000
Di-n-octylphthalate	44000 UD	44000
Benzo(b) fluoranthene	22000 JD	44000
Benzo(k)fluoranthene	7700 JD	44000
Benzo(a)pyrene	15000 JD	44000
Indeno(1,2,3-cd)pyrene	14000 JD	44000
Dibenz(a,h)anthracene	44000 UD	44000
Benzo(g,h,i)perylene	14000 JD	44000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	40%J D	25-121%
Phenol-d5	69 % JD	24-113%
2-Chlorophenol-d4	61 % JD	20-130%
1,2-Dichlorobenzene-d4	76 % JD	20-130%
Nitrobenzene-d5	156%JD	23-120%
2-Fluorobiphenyl	9 2% JD	30-115%
2,4,6-Tribromophenol	88 % JD	19-122%
Terphenyl-d14	132%JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLMO1)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

Huntingdon_

HPN: 2757



ecology and environment, inc.

111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-663-9415
International Specialists in the Environment

MEMORANDUM

DATE: March 24, 1994

TO: John Nordine, Project Manager, E & E, Chicago, IL

•

FROM: David Hendren, TAT-Chemist, E & E, Chicago, IL X

THRU: Pat Zwilling, ATATL, E & E, Chicago, IL

SUBJ: Polychlorinateddibenzo Dioxins and Furans Data Quality

Assurance Review for Scrap Metal site, Chicago, Cook

County, Illinois

REF: Analytical TDD:T05-9402-807 Project TDD:T05-9402-007

Analytical PAN: EIL0831AAA Project PAN: EIL0831SAA

The data quality assurance review of 4 soil samples collected from the site has been completed. Analysis for polychlorinated dibenzodioxins (PCDD) and polychlorinated dibenzofurans (PCDF) was performed by Twin City Testing Corporation (Huntingdon) of St. Paul, Minnesota following U.S. EPA SW-846 Method 8290.

The samples were numbered as follows in the field. The corresponding laboratory identification number is provided:

TAT Sample #	corresponds	to	-	>	Laboratory Sa	ample	土
S1					15270		
S2					15277		
S3					15278		
S8					15290		

Data Qualifications:

I. Holding Time: Acceptable

The samples were collected on 2/22/94, extracted on 3/4/94 and analyzed on 3/7/94 and 3/9/94. The holding time criteria of 6 months from sample collection to extraction was satisfied. All

samples were analyzed within the 40 day limit following extraction.

II. Instrument Performance: Acceptable

Analysis of the calibration check solution showed that the % valley between tetrachlorodibenzodioxin (TCDD) isomers (1,2,3,4-) and (2,3,7,8-) was less than 25 %, as required.

III. Calibration:

A. Initial Calibration: Acceptable

The percent relative standard deviations (%RSD) for all analytes in the initial calibration (performed 11/9/93) was less than 15%, as required, except for OCDD (15.66%). Qualification was not judged to be necessary. All specific ion ratio criteria were achieved and signal to noise criteria were satisfied.

B. Continuing Calibration: Acceptable

The continuing calibration standard was analyzed before and following sample analysis. All percent difference of response factors were less than 30%, except for OCDD(34.1%) in the final calibration check run on 3/7/94. Since the initial calibration check was in control and all other parameters were acceptable, qualification was not judged to be necessary.

IV. Method Blank: Acceptable

A method blank was analyzed for each matrix analyzed. HxCDF, HpCDD, and OCDD were detected at 0.39 PPT, 4.4 PPT and 21.0 PPT respectively. All reported values in the samples for these analytes exceeded these levels (over five times) and qualification was not required. No other target analytes or contaminants were detected above the detection limit.

V. Internal Standards: Acceptable

The recoveries of the 13C-internal standards were acceptable in all samples.

VI. Matrix Spike/Matrix Spike Duplicate (MS/MSD): Acceptable

The percent recoveries for the Matrix Spike/Matrix Spike Duplicate (MS/MSD) were within the established quality control limits.

VII. Overall Assessment of Data For Use:

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality Control Guidance for Removal Activities" (OSWER 9360.4-01 April, 1990). Based upon the information provided, the data are acceptable for use.

******** *TWIN CITY TESTING CORPORATION* *METHOD 8290 ANALYSIS RESULTS * ***** Client....ECOLOGY & ENVIRONMENT

Client's Sample ID.....S1 (1:2.5 DILUTION)
TCT Sample ID.......15270
Analysis Date......3/7/94 12:35
Filename......V40307D
Injected By.......MJK Total Amount Extracted...0.0129 kg Method Blank ID.....BLANK-15828 Extraction Date.....3/4/94

NATIVE ISOMERS	CONC. ng/kg	LOD ng/kg	INTERNAL STANDARDS	ng's ADDED	PERCENT RECOVERY
2378-TCDF TOTAL TCDF	220 * 1500		2378-TCDF-13C 2378-TCDD-13C 12378-PeCDF-13C	2.00 2.00 2.00	105 98 87
2378-TCDD TOTAL TCDD	ND 34	11	23478-PeCDF-13C 12378-PeCDD-13C 123478-HxCDF-13C.	2.00 2.00 2.00	91 81 102
12378-PeCDF 23478-PeCDF TOTAL PeCDF	94 260 1300		123678-HXCDF-13C. 234678-HXCDF-13C. 123789-HXCDF-13C. 123478-HXCDD-13C.	2.00 2.00 2.00 2.00	81 93 81 98
12378-PeCDD TOTAL PeCDD	ND ND	310	123478-HXCDD-13C. 1234678-HPCDF-13C 1234789-HPCDF-13C	2.00 2.00 2.00	97 73 72
123478-HxCDF 123678-HxCDF 234678-HxCDF	250 160 230		1234678-HPCDD-13C OCDD-13C	2.00 4.00	82 82
123789-HxCDF TOTAL HxCDF	95 2000		1234-TCDD-13C 123789-HxCDD-13C.	2.00 2.00	NA NA
123478-HxCDD 123678-HxCDD 123789-HxCDD	21 70 33		2378-TCDD-37C14	0.20	INT
TOTAL HXCDD	240	~~~~	Total 2378-TCDD Equivalence:	265_	ng/kg
1234678-HpCDF 1234789-HpCDF TOTAL HpCDF	900 120 1500		(Using ITE Facto	ors/DB-5	Data)
1234678-HpCDD TOTAL HpCDD	730 1400				
OCDF	770 43 00				

* Value may include contributions from other TCDF isomers.

All values are expressed on a dry weight basis.

CONC= Concentration (Totals include 2378-substituted isomers.) LOD = Limit of Detection

ND = Not Detected
NA = Not Applicable
INT = Interference

TCT Invoice Number....4411 94-2757



****************** *TWIN CITY TESTING CORPORATION* *METHOD 8290 ANALYSIS RESULTS * Client....ECOLOGY & ENVIRONMENT

Client's Sample ID.....S2 (1:20 DILUTION)
TCT Sample ID......15277
Analysis Date......3/9/94 14:42

CCAL Filename........V40309B

Method Blank ID.....BLANK-15828

Extraction Date.....3/4/94

NATIVE ISOMERS	CONC. ng/kg	LOD ng/kg	INTERNAL STANDARDS	ng's ADDED	PERCENT RECOVERY
2378-TCDF TOTAL TCDF	ND 17000	10000	2378-TCDF-13C 2378-TCDD-13C 12378-PeCDF-13C	2.00 2.00 2.00	103 60 89
2378-TCDD TOTAL TCDD	ND 580	22	23478-PeCDF-13C 12378-PeCDD-13C 123478-HxCDF-13C.	2.00 2.00 2.00	120 97 128
12378-PeCDF 23478-PeCDF TOTAL PeCDF	4400 6300 42000		123478-HXCDF-13C. 123678-HXCDF-13C. 234678-HXCDF-13C. 123789-HXCDF-13C. 123478-HXCDD-13C.	2.00 2.00 2.00 2.00 2.00	88 94 98 98
12378-PeCDD TOTAL PeCDD	ND ND	4200	123678-HxCDD-13C. 1234678-HpCDF-13C 1234789-HpCDF-13C	2.00 2.00 2.00	90 95 71
123478-HxCDF 123678-HxCDF 234678-HxCDF	ND ND 2700	20000 4100	1234678-HPCDD-13C OCDD-13C	2.00 4.00	72 47
123789-HXCDF TOTAL HXCDF	890 15000		1234-TCDD-13C 123789-HxCDD-13C.	2.00	NA NA
123478-HxCDD 123678-HxCDD 123789-HxCDD	ND 250 130	140	2378-TCDD-37C14	0.20	INT
TOTAL HXCDD	380		Total 2378-TCDD Equivalence:	4004	ng/kg
1234678-HpCDF 1234789-HpCDF TOTAL HpCDF	12000 3600 16000		(Using ITE Fact	ors/DB-5	Data)
1234678-HpCDD TOTAL HpCDD	2900 5900				
OCDF OCDD	30000 220 00				

All values are expressed on a dry weight basis.

CONC= Concentration (Totals include 2378-substituted isomers.)
LOD = Limit of Detection

ND = Not Detected NA = Not Applicable

TCT Invoice Number....4411 94-2757 INT = Interference



```
*TWIN CITY TESTING CORPORATION*
                      *METHOD 8290 ANALYSIS RESULTS *
                       Client....ECOLOGY & ENVIRONMENT
Client's Sample ID.....S3 (1:2.5 DILUTION)
TCT Sample ID........15278
Analysis Date........3/7/94_14:30
Total Amount Extracted...0.0155 kg
Method Blank ID.....BLANK-15828
Extraction Date.....3/4/94
                                            INTERNAL
                                                            ng's
                                                                    PERCENT
NATIVE
                CONC.
                            LOD
                                                           ADDED
                                                                    RECOVERY
ISOMERS
                ng/kg
                           ng/kg
                                            STANDARDS
2378-TCDF
                  1700 *
                                       2378-TCDF-13C....
                                                             2.00
                                                                        121
                                       2378-TCDD-13C....
12378-PeCDF-13C...
23478-PeCDF-13C..
                                                             2.00
TOTAL TCDF
                                                                        111
                  9400
                                                             2.00
                                                                        116
                                                             2.00
                                                                        124
                              18
2378-TCDD
                    ND
                                       12378-PeCDD-13C..
                                                             2.00
TOTAL TCDD
                   130
                                                                        106
                                       123478-HxCDF-13C.
123678-HxCDF-13C.
                                                                        120
                                                             2.00
                                                                         91
                  550
                                                             2.00
12378-PeCDF
                                       234678-HxCDF-13C.
                                                             2.00
23478-PeCDF
                  1300
                                                                        114
                                       123789-HxCDF-13C.
                                                             2.00
                                                                         95
TOTAL PeCDF
                  7600
                                       123478-HxCDD-13C.
                                                             2.00
                                                                        114
                                       123678-HxCDD-13C.
                                                             2.00
                                                                         99
12378-PeCDD
                    ND
                             160
                                       1234678-HpCDF-13C
                                                             2.00
                                                                         87
TOTAL PeCDD
                    ND
                                       1234789-HpCDF-13C
                                                             2.00
                                                                         84
                                       1234678-HpCDD-13C
                                                             2.00
                                                                         87
123478-HxCDF
                   810
                                       OCDD-13C.....
                                                             4.00
                                                                         78
123678-HxCDF
                   640
                   990
220
234678-HxCDF
                                                             2.00
                                                                         NA
                                       1234-TCDD-13C....
123789-HxCDF
                  7700
                                       123789-HxCDD-13C.
                                                             2.00
                                                                         NA
TOTAL HXCDF
                                       2378-TCDD-37C14..
                                                             0.20
                                                                        INT
                   100
123478-HxCDD
                           ----- ,
123678-HxCDD
                   170
123789-HxCDD
                   120
                                          Total 2378-TCDD
                  1100
TOTAL HXCDD
                                          Equivalence:
                                                             1207
                                                                   ng/kg
                                          (Using ITE Factors/DB-5 Data)
                  3100
1234678-HpCDF
                           ____
1234789-HpCDF
                   330
TOTAL HpCDF
                  4500
1234678-HpCDD
                  1500
TOTAL HpCDD
                  2900
OCDF
                  1600
OCDD
                  3900
* Value may include contributions from other TCDF isomers.
All values are expressed on a dry weight basis.
      Concentration (Totals include 2378-substituted isomers.)
CONC=
LOD = Limit of Detection
ND = Not Detected
NA = Not Applicable
```



INT = Interference

TCT Invoice Number....4411 94-2757

TWIN CITY TESTING CORPORATION *METHOD 8290 ANALYSIS RESULTS * Client....ECOLOGY & ENVIRONMENT

Client's Sample ID......S8 (1:20 DILUTION) TCT Sample ID.........15290

Injected By........MJK
Total Amount Extracted...0.0144 kg % Moisture..... 29.2 % ICAL Date......11/9/93 CCAL Filename......V40309B
Method Blank ID.....BLANK-15828

Extraction Date.....3/4/94

NATIVE ISOMERS	conc. ng/kg	LOD ng/kg	INTERNAL STANDARDS	ng's ADDED	PERCENT RECOVERY
2378-TCDF TOTAL TCDF	ND 1900	380	2378-TCDF-13C 2378-TCDD-13C 12378-PeCDF-13C	2.00 2.00 2.00	87 79 99
2378-TCDD TOTAL TCDD	ND 150	17	23478-PeCDF-13C 12378-PeCDD-13C 123478-HxCDF-13C.	2.00	132 93 105
12378-PeCDF 23478-PeCDF TOTAL PeCDF	ND 220 3300	6000	123478-HXCDF-13C. 123678-HXCDF-13C. 234678-HXCDF-13C. 123789-HXCDF-13C. 123478-HXCDD-13C.	2.00 2.00 2.00 2.00	75 91 87 92
12378-PeCDD TOTAL PeCDD	ND ND	31	123678-HxCDD-13C. 1234678-HpCDF-13C 1234789-HpCDF-13C	2.00 2.00 2.00	86 83 67
123478-HxCDF 123678-HxCDF 234678-HxCDF	500 ND 360	340	1234678-HPCDD-13C OCDD-13C	2.00 4.00	77 60
123789-HxCDF TOTAL HxCDF	170 2700		1234-TCDD-13C 123789-HxCDD-13C.	2.00 2.00	NA NA
123478-HxCDD 123678-HxCDD 123789-HxCDD	29 670 190		2378-TCDD-37C14	0.20	INT
TOTAL HXCDD	1800 1100		Total 2378-TCDD Equivalence: (Using ITE Fact	360	ng/kg Data)
1234678-HpCDF 1234789-HpCDF TOTAL HpCDF	240 1300		(USING TIE FACE	015/06-3	Dacay
1234678-HpCDD TOTAL HpCDD	2800 5700				
OCDF OCDD	1600 15000				

All values are expressed on a dry weight basis.

CONC= Concentration (Totals include 2378-substituted isomers.)

LOD = Limit of Detection

ND = Not Detected

NA = Not Applicable INT = Interference TCT Invoice Number....4411 94-2757



APPENDIX C RCMS COST ESTIMATE

(Portions of this appendix were redacted which were not relevant to the Removal Action at this site.)

Page: 1

Summary Report Initial Cost Projection Scenario: SCRAP METAL

Projection ID Number: IL0831SA Date: 05/02/94

Cleanup Contractor: ESE5 - ESE TAT Contractor: ECOLOGY&ENVIRON

Cost Projection Summary -----

Contractor Personnel	282,065.90
Contractor Equipment	55,316.98
Unit Rate Materials	115,944.00
At Cost Materials	1,302.50
Subcontractors	14,220.00
Waste Transportation	7,312.50
Waste Disposal	702,000.0 0
Cleanup Contractor Subtotal	1,178,161.88
Federal and State Agencies	0.00
Extramural Subtotal	1,178,161.88
15 % Extramural Contingency	176,724.28
Extramural Subtotal	1,354,886.16
TAT Personnel	97,632.00
TAT Special Projects	0.00
TAT Analytical Services	0.00
Total TAT Costs	97,632.00
Other Cost Items	0.00
Evenement Cubectal	
Extramural Subtotal	1,452,518.16
10 % Project Contingency	145,251.82
Total Extramural Cost	1,597,769.98
EPA Regional Personnel	32,400.00
EPA Non-Regional Personnel	0.00
EPA Headquarters Direct	0.00
(0 % of Regional Hours)	
EPA Indirect	65,880.00
EPA Total	98,280.00
Project Total	1,696,049.98

Page: 2

Summary Report (cont.) Initial Cost Projection Scenario: SCRAP METAL

Projection ID Number: IL0831SA Date: 05/02/94

Cleanup Contractor: ESE5 - ESE TAT Contractor: ECOLOGY&ENVIRON

Project Scope

		Estimated	
Number	Step/Milestone	Duration	Cost
000	GENERAL SITE COSTS	120 Days	1,696,049.98
999	ARCHIVE COSTS	120 Days	0.00
	: :		1,696,049.98